

## **SLURM Administrators Tutorial**

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Resource Management Systems Architect

- Introduction
- SLURM scalable and flexible RJMS
- Part 1: Basics
  - Overview, Architecture, Configuration files, Partitions, Plugins, Reservations, Reconfiguration
- Part 2: Advanced Configuration
  - Accounting, Scheduling, Allocation, Network Topology Placement, Generic Resources, Licenses Management, Energy Reduction Techniques
- Part 3: Experts Configuration
  - CPU Management, Isolation with cgroups, Power Management, Simulation and evaluation
- Upcoming Features

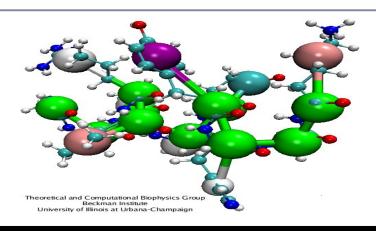


#### Introduction

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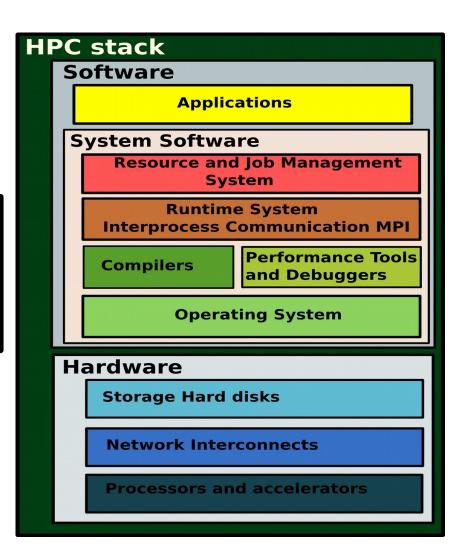
# **High Performance Computing**



## **System Software:**

Operating System, Runtime System, Resource Management, I/O System, Interfacing to External Environments





## **Resource and Job Management Systems**

The goal of a Resource and Job Management System (RJMS) is to satisfy users' demands for computation and assign resources to user jobs with an efficient manner.

# Resource and Job Management System Hardware Nodes

## **RJMS Importance**

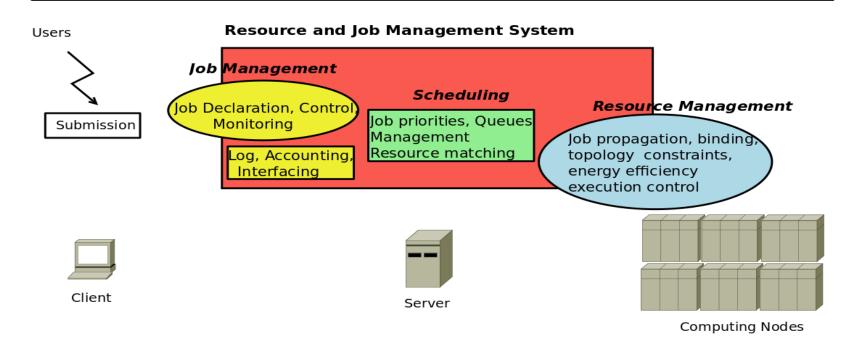
Strategic Position but complex internals:

- Direct and constant knowledge of resources
- Multifacet procedures with complex internal functions

## Resource and Job Management System Layers

This assignement involves three principal abstraction layers:

- •Job Management: declaration of a job and demand of resources and job characteristics,
- •Scheduling: matching of the jobs upon the resources,
- •Resource Management: launching and placement of job instances upon the computation resources along with the job's control of execution



# **Resource and Job Management System Concepts**

RJMS subsystems	Principal Concepts	Advanced Features
Resource Management	-Resource Treatment (hierarchy, partitions,) -Job Launcing, Propagation, Execution control -Task Placement (topology,binding,)	- High Availability - Energy Efficiency - Topology aware placement
Job Management	-Job declaration (types, characteristics,) -Job Control (signaling, reprioritizing,) -Monitoring (reporting, visualization,)	- Authentication (limitations, security,) - QOS (checkpoint, suspend, accounting,) - Interfacing (MPI libraries, debuggers, APIs,)
Scheduling	-Scheduling Algorithms (builtin, external,) -Queues Management (priorities,multiple,)	- Advanced Reservation

#### Introduction

#### SLURM scalable and flexible RJMS

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## **SLURM scalable and flexible RJMS**

- •SLURM **open-source** Resource and Job Management System, sources freely available under the GNU General Public License.
- •Portable: written in C with a GNU autoconf configuration engine.
- •Modular: Based on a plugin mechanism used to support different kind of scheduling policies, interconnects, libraries, etc
- •Robust: highly tolerant of system failures, including failure of the node executing its control functions.
- •Scalable: designed to operate in a heterogeneous cluster with up to tens of millions of processors. It can accept 1,000 job submissions per second and fully execute 500 simple jobs per second (depending upon hardware and system configuration).
- •Power Management: Job can specify their desired CPU frequency and power use by job is recorded. Idle resources can be powered down until needed.

## **SLURM History and Facts**

- •Initially developed in LLNL since 2003, passed to SchedMD in 2011
- •Multiple enterprises and research centers have been contributing to the project (LANL, CEA, HP, BULL, BSC, CRAY etc)
- •Large international community, active mailing lists (support by main developers)
  - Contributions (various external software and standards are integrated upon SLURM)
- As of the June 2014 Top500 supercomputer list, SLURM is being used on six of the ten most powerful computers in the world including the no1 system, Tianhe-2 with 3,120,000 computing cores.





#### **BULL and SLURM**

- BULL initially started to work with SLURM in 2005
- About 6 SLURM-dedicated engineers since 2013
  - Research upon the field of Resource Management and Job Scheduling (National/European financed projects, PhDs) and definition of RoadMap
  - Development of new SLURM features: all code dropped in the open-source
  - Support upon clusters : Training, Confgiruation, Bugs, Feature Requests, etc
- Integrated as the default RJMS into the BULL- HPC software stack since 2006
- Close development collaboration with SchedMD and CEA
- Organaziation of Slurm User Group (SUG) Conference (User, Admin Tutorials + Technical presentation for developpers)
   http://www.schedmd.com/slurmdocs/publications.html

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# **SLURM** sources and Documentation

#### **Slurm sources:**

- Download a repo (stable or development) from: http://www.schedmd.com/#repos
- Or the latest code from: git clone git://github.com/SchedMD/slurm.git

-For User and Admins latest **documentation**:

http://slurm.schedmd.com/documentation.html

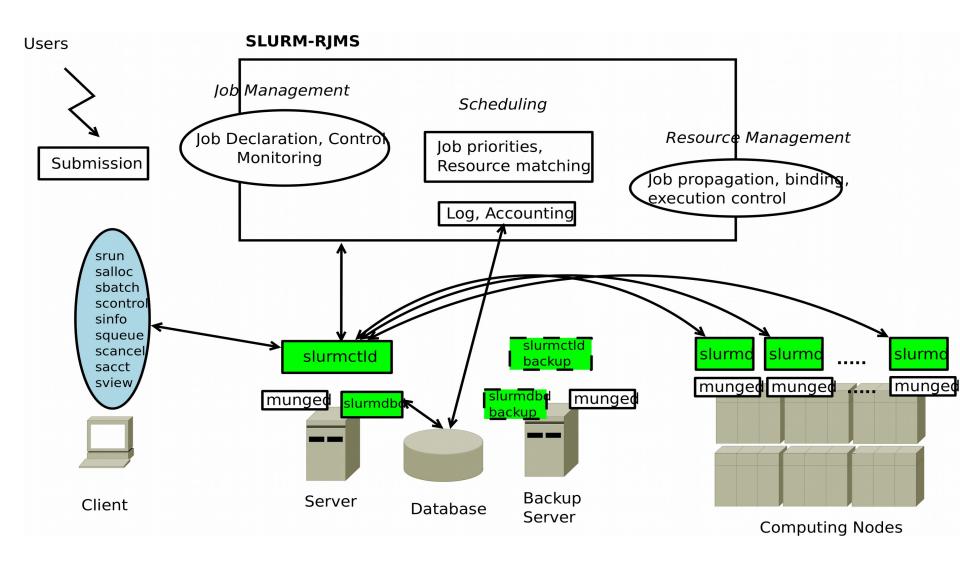
- -Detailed **man pages** for commands and configuration files http://slurm.schedmd.com/man\_index.html
- -All SLURM related **publications and presentations**:

http://slurm.schedmd.com/publications.html

# **Quick Installation Steps**

- 1.Install MUNGE for authentication. Make sure the MUNGE daemon, munged is started before you start the SLURM daemons.
- 2.Install SLURM either creating a tgz from git or downloading an existing tgz
- 3.cd to the directory containing the SLURM source and type ./configure with necessary options.
- 4. Type *make* to compile SLURM.
- 5. Type *make install* to install the programs, documentation, libraries, header files, etc.
- 6.Create the slurm User upon all nodes of the cluster.
- 7.Create parent directories for SLURM's log files, process ID files, state save directories, etc. are not created by SLURM. They must be created and made writable by SlurmUser as needed prior to starting SLURM daemons.
- 8. Create a basic slurm.conf file
- 9. Follow the presentation...

#### **SLURM Architecture**



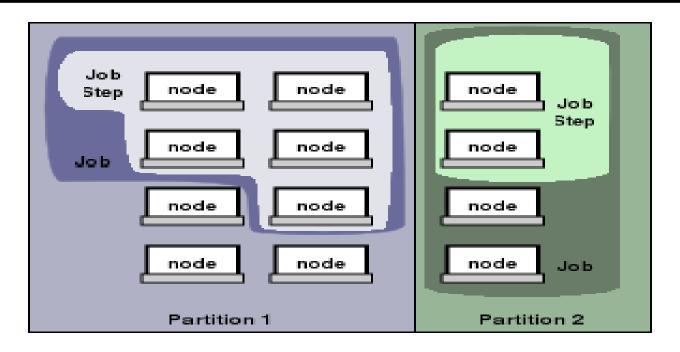
#### **SLURM Terms**

Computing node Computer used for the execution of programs

Partition Group of nodes into logical sets

•Job allocation of resources assigned to a user for some time

•Step sets of (possible parallel) tasks with a job



# **SLURM Principles**

## Architecture Design:

- one central controller daemon slurmctld
- A daemon upon each computing node slurmd
- One central daemon for the database controls slurmdbd

## Principal Concepts:

- a general purpose plugin mechanism (for features such as scheduling policies, process tracking, etc)
- the partitions which represent group of nodes with specific characteristics (job limits, access controls, etc)
- one queue of pending work
- The job steps which are sets of (possibly parallel) tasks within a job

#### **User Commands**

- **srun** allocate resources ( number of nodes, tasks, partition, constraints, etc.) launch a job that will execute on each allocated cpu.
- salloc allocate resources (nodes, tasks, partition, etc.), either run a command or start a shell. Request launch srun from shell. (interactive commands within one allocation)
- **sbatch** allocate resources (nodes, tasks, partition, etc.) Launch a script containing sruns for series of steps.
- **sbcast** transmit file to all nodes of a running job. Used in sbatch or salloc.

**sattach** attach to running job for debuggers.

#### **User & Admin Commands**

```
sinfo display characteristics of partitions
squeue display jobs and their state
scancel cancel a job or set of jobs.
scontrol display and changes characteristics of jobs, nodes, partitions.
sstat show status of running jobs.
sacct display accounting information on jobs.
sprio show factors that comprise a jobs scheduling priority
smap graphically show information on jobs, nodes, partitions
```

## **Admin Commands**

- **sacctmgr** setup accounts, specify limitations on users and groups.
- **sreport** display information from accounting database on jobs, users, clusters.
- **sview** graphical view of cluster. Display and change characteristics of jobs, nodes, partitions.
- **strigger** show, set, clear event triggers. Events are usually system events such as an equipement failure.
- **sshare** view sharing information from multifactor plugin.

# **Configuration**

Slurm configuration: Through configuration files responsible, for the function of different daemons present on the management and the computing nodes

#### slurm.conf

• Indispensable on all nodes (management-compute)

#### slurmdbd.conf

- Used if slurmdbd accounting
- Only on management node

## topology.conf

- Used if topology plugin activated
- Indispensable on all nodes (management-compute)

## gres.conf

- Used if gres plugin activated
- Only on computing nodes

## cgroup.conf

- Used if cgroup plugin activated
- Only on computing nodes

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#### slurm.conf

- Low level configuration
- Management policies
- Scheduling policies
- Allocation policies
- Node definition
- Partition definition

#### slurmdbd.conf

- Type of persistent storage (DB)
- Location of storage

## topology.conf

Switch hierarchy

#### gres.conf

- Generic resources details
- Device files

## cgroup.conf

- Mount point
- Release agent path
- Cgroup subsystems parameters

	Controller	Compute node
Mandatory	slurm.conf slurmdbd.conf	slurm.conf
Optional	prologs epilogs topology.conf	gres.conf cgroup.conf topology.conf

# Configuration (slurm.conf) - Part 1

#### **Node definition**

- Characteristics (sockets, cores, threads, memory, features)
- Network addresses

#### **Partition definition**

- Set of nodes
- Sharing
- Priority/preemption

#### # Compute Nodes

NodeName=cuzco[1-10] Procs=16 Sockets=2 CoresPerSocket=8 ThreadsPerCore=1 State=UNKNOWN RealMemory=38000 NodeName=cuzco[10-20] Procs=32 Sockets=2 CoresPerSocket=8 ThreadsPerCore=2 State=UNKNOWN RealMemory=46000

#### # Partitioning

PartitionName=exclusive Nodes=cuzco[1-20] MaxTime=INFINITE State=UP Priority=10 Shared=Exclusive PartitionName=shared Nodes=berlin[1-20] Default=YES MaxTime=INFINITE State=UP Priority=30 PartitionName=procs16 Nodes=berlin[1-10] MaxTime=INFINITE State=UP Priority=30 PartitionName=procs32 Nodes=berlin[10-20] MaxTime=INFINITE State=UP Priority=30

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## **Partitions**

Partitions are used in SLURM to group nodes/resources characteristics

**Partition 1:** 32 cores and high\_memory

Partition 2: 32 cores and low\_memory

Partition 3: 64 cores

## **More on Partitions**

## **Shared Option**

Controls the ability of the partition to execute more than one job on a resource (node, socket, core)

**EXCLUSIVE** allocates entire node (overrides cons\_res ability to allocate cores and sockets to multiple jobs)

**NO** sharing of any resource.

**YES** all resources can be shared, unless user specifies –exclusive on srun | salloc | sbatch

**Important Note:** To view the particular parameters of partitions users can use the "scontrol show partitions" command

## **Configuration (slurm.conf) - Part 2**

#### #slurm.conf

#### # Basic parameters

ClusterName=cuzco
ControlMachine=cuzco0
#ControlAddr=127.0.0.1
SlurmUser=slurm
SlurmctldPort=6817
SlurmdPort=6818
AuthType=auth/munge

#### # States saving

StateSaveLocation=/var/spool/slurm
SlurmdSpoolDir=/var/spool/slurmd.%n
SlurmctldPidFile=/var/run/slurmctld.pid
SlurmdPidFile=/var/run/slurmd.%n.pid

#### # Logging

SlurmctldDebug=5
SlurmctldLogFile=/var/log/slurmctld.log
SlurmdDebug=5
SlurmdLogFile=/var/log/slurmd.%n.log

#### # Timers

SlurmctldTimeout=300
SlurmdTimeout=300

## **Management Policies**

- Location of controllers, spool, state info
- Authentication
- Logging
- Prolog / epilog scripts

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# **Configuration (slurm.conf) - Part 3**

#### # Process-Task tracking

ProctrackType=proctrack/linuxproc TaskPlugin=task/affinity TaskPluginParam=Cpusets

#### **# Selection of Resources**

SelectType=select/cons\_res
SelectTypeParameters= CR\_Core\_Memory

#### # Scheduling

SchedulerType=sched/backfill FastSchedule=1 PreemptMode=REQUEUE PreemptType=preempt/qos FastSchedule=1

## **Scheduling policies**

- Priority
- Preemption
- Backfill

## **Allocation policies**

- Entire nodes or 'consumable resources'
- Task Affinity (lock task on CPU)
- Topology (minimum number of switches)

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# **Plugins in SLURM**

- -Authentication (i.e. munge,)
- -Job Accounting Gather (i.e. linux, cgroups)
- -Accounting Storage (i.e. mysql, postgres)
- -Generic Resources (GRES) (i.e. gpu, nic)
- -Job Submission (i.e. partitions, lua)
- -MPI (i.e. openmpi, pmi2)
- -Energy Accounting (i.e. rapl,ipmi)
- -Preemption (i.e. partitions,qos)
- -Priority (i.e. basic, multifactor)
- -Process Tracking (i.e. linux,cgroup)
- -Scheduler (i.e. builtin,backfill)
- -Resource Selection (i.e. linear,cons\_res)
- **-Task** (i.e. affinity,cgroups)
- -Topology (i.e. tree,3d\_torus)

## **Examples of info commands**

```
> sinfo
PARTITION AVAIL
                TIMELIMIT
                           NODES STATE NODELIST
all*
            up infinite
                               4 idle trek[0-3]
P2
            up infinite 4 idle trek[0-3]
                               4 idle trek[0-3]
P3
            up infinite
> scontrol show node trek
NodeName=trek3 Arch=x86 64 CoresPerSocket=4
   CPUAlloc=0 CPUErr=0 CPUTot=16 Features=HyperThread
  Gres=(null)
   NodeAddr=sulu NodeHostName=sulu
   OS=Linux RealMemory=1 Sockets=2
   State=IDLE ThreadsPerCore=2 TmpDisk=0 Weight=1
   BootTime=2011-06-30T11:04:22 SlurmdStartTime=2011-07-12T06:23:43
   Reason=(null)
> scontrol show partition P2
PartitionName=P2
  AllocNodes=ALL AllowGroups=ALL Default=NO
   DefaultTime=NONE DisableRootJobs=NO GraceTime=0 Hidden=NO
   MaxNodes=UNLTMTTFD MaxTime=UNLTMTTFD MinNodes=1
   Nodes=trek[0-3]
  Priority=2 RootOnly=NO Shared=NO PreemptMode=CANCEL
   State=UP TotalCPUs=40 TotalNodes=4
```

## srun, salloc, sbatch

- All request an allocation of resources.
- Similar set of command line options.
- Request number of nodes, tasks, cpus, constraints, user info, dependencies, and lots more.
- srun launches tasks (command) on the requested nodes.
- salloc can launch a task such as mpirun on the client, or open a shell on the client. Srun is then used to launch tasks within the allocation.
- sbatch is a shell script that contains multiple sruns within the allocation. (multi step job).

## **Example srun**

# srun -l -p P2 -N2 -tasks-per-node=2 -exclusive hostname

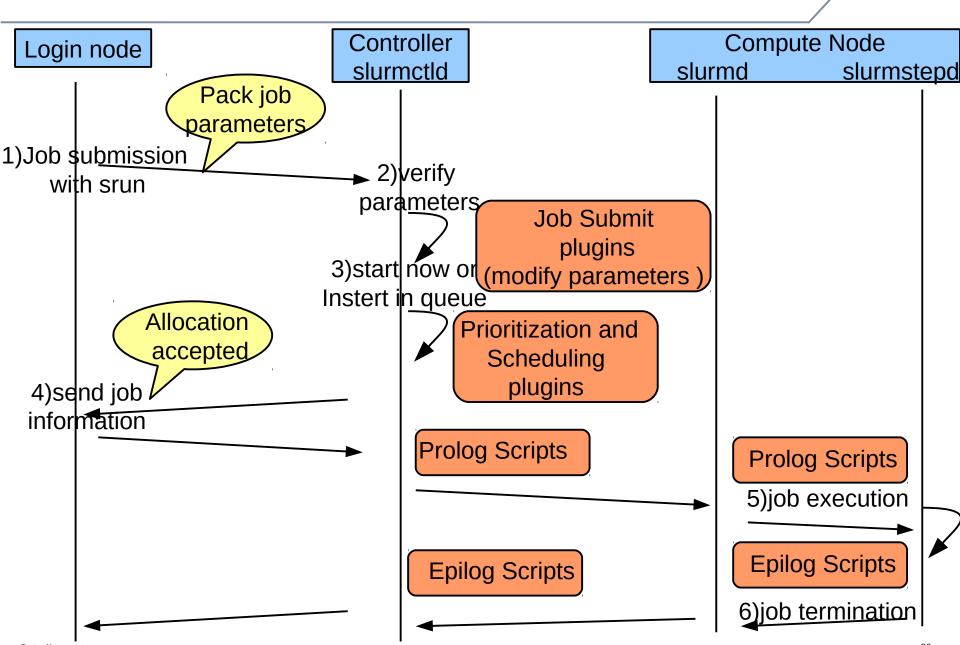
- -l prepend task number to output (debug)
- -p P2 use Partition P2
- -N2 use 2 nodes
- --tasks-per-node launch 2 tasks on each node
- --exclusive do not share the nodes

hostname command to run.

#### Results

- > srun -1 -p P2 -N2 --tasks-per-node=2
- -- exclusive hostname
- 0: scotty
- 1: scotty
- 2: bones
- 3: bones

# **Job Journey**



#### login node :

Using srun, all information about the job (parameters) are packed and send via RPC to the controller

#### Slurm controller :

- \_ call job\_submit plugin to modify parameters (if needed)
- \_ validate all the parameters (partition, limits, etc...)
- \_ evaluate the priority of the job: it's the highest > start now,

it's not > enqueue

Direct path and « queued » path :

- \_ find nodes matching job's recquirement
- \_ give credential to srun → job can start
- \_ update nodes state

#### Job journey - we have compute nodes, what's next?

- srun sends all job's information (executable name, files name, etc...) to the first node of the allocation.
- Then the node forwards the request to other nodes using a spanning tree algorithm
- Each slurmd send a request to the controller to verify the job was not canceled during the process
- Each slurmd spawn a slurmstepd program to manage the job step
- Slurmstepd can now initiate the job :
  - do the mpi setup
  - \_ create a container using the proctrack plugin
  - \_ change the uid to the user one
  - \_ set up environnement variables
  - fork the task

#### Job journey - This is the end...

- job termination can be initiated by :
  - the task itself : end of the program
    - → Slurmd will warn slurmctld which will log the termination
  - the user: job cancelation,
  - Slurm: the job has exceed a limit (timelimit, memory usage, etc...)
    - $\rightarrow$  in both case slurmd will:
      - \_ send a SIGTERM to any task
      - \_ wait kilwait seconds
      - \_ send a SIGKILL
      - \_ execute any Epilog
      - \_ send a final message to slurmctld

# **JobSubmit Plugin in SLURM**

## JobSubmitPlugins Parameter in slurm.conf

A comma delimited list of job submission plugins to be used. These are intended to be site-specific plugins which can be used *to* set default job parameters and/or logging events. Plugins that can be modified and used: "defaults", "partition", "lua",etc. For examples of use, see the SLURM code in "src/plugins/job\_submit" and "contribs/lua/job\_submit\*.lua" then modify the code to satisfy your needs.

#### Example:

Associate user jobs to projects: A user may only charge jobs to one of his projects

Parameter	Location	Invoked by	User	When executed
Prolog (from slurm.conf)	Compute or front end node	slurmd daemon	SlurmdUser (normally user root)	First job or job step initiation on that node (by default);
PrologSlurmctl d (from slurm.conf)	Head node (where slurmctld daemon runs)	slurmctld daemon	SlurmctldUser	At job allocation
Epilog (from slurm.conf)	Compute or front end node	slurmd daemon	SlurmdUser (normally user root)	At job termination
EpilogSlurmctld (from slurm.conf)	Head node (where slurmctld daemon runs)	slurmctld daemon	SlurmctldUser	At job termination

Parameter	Location	Invoked by	User	When executed
SrunProlog (from slurm.conf) or srunprolog	srun invocation node	srun command	User invoking srun command	Prior to launching job step
TaskProlog (from slurm.conf) or srun –task- prolog	Compute node	slurmstepd daemon	User invoking srun command	Prior to launching job step
TaskEpilog (from slurm.conf) or srun –task- epilog	Compute node	slurmstepd daemon	User invoking srun command	Completion job step
SrunEpilog (from slurm.conf) or srunepilog	srun invocation node	srun command	User invoking srun command	Completion job step

- Once the principal configuration parameters are correctly set the services can be started on management and computing nodes by launching the particular scripts on all nodes: /etc/init.d/slurm {start, stop, restart, ...}
- •Alternatively the services can be started by **executing the commands** slurmctld on the controller and slurmd on the computing nodes
- •The services are normally launched in the background with logging in the particular files set in the slurm.conf. However it is possible to **start the deamons in the foreground** with -D followed by v for different verbosity levels. This is useful for testing.

slurmctld -Dvvvvvv slurmd -Dvvvvvv

- Every **change of** configuration parameters should be followed by a **copy of slurm.conf** on all computing nodes and restart of SLURM deamons.
- In case other configuration files are changed they should be equally copied on all computing nodes (except of slurmdbd.conf which is used only on the management nodes)
- Reconfiguration can also be enabled dynamically (running jobs continue execution) without copying the slurm.conf file with scontrol reconfigure (doesn't update slurm.conf files)
- ALL SLURM daemons should be shutdown and restarted if any of these parameters are to be changed: AuthType, BackupAddr, BackupController, ControlAddr, ControlMach, PluginDir,StateSaveLocation, SlurmctldPort or SlurmdPort

### scontrol

command can be also used for reservations
It provides the ability to create, update and delete
advanced reservations for resources allocations

Basic parameters that need to be used:

Starttime, Duration, User, NodeCnt or NodeList

Once the reservation is made the user can submit a job upon the reserved Resources and this job will start on the starttime.

### **Examples:**

>scontrol: create res StartTime=2009-04-01T08:00:00 Duration=5:00:00 Users=toto NodeCnt=10 Reservation created: toto 1

>scontrol: update Reservation=toto\_1 Flags=Overlap NodeCnt=20

An alternative way to start a job in a particular moment in the future is the –begin-time option of the submission commands

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# **SLURM Accounting**

- Accounting based upon Mysql database
- Robust and scalable (confirmed upon Tera100 cluster)
- Command for database and accounting configuration: sacctmgr
- Fairsharing and Preemption scheduling techniques based upon the accounting infrastructure

```
>sacctmgr add cluster snowflake
>sacctmgr add account users Cluster=snowflake Description="none" Organization="none"
>sacctmgr list users
User Def Acct Admin

gohn students None
root root Administ+
slurm professors None
```

# **Accounting**

### **Commands**

**Sacct** reports resource usage for running or terminated jobs.

**Sstat** reports on running jobs, including imbalance between tasks.

**Sreport** generates reports based on jobs executed in a time interval.

Sacctmgr is used to create account and modify account settings.

## Plugins associated with resource accounting

**AccountingStorageType** controls how information is recorded (MySQLI with SlurmDBD is best)

**JobAccntGatherType** controls the mechanism used to gather data. (OS Dependent)

**JobCompType** controls how job completion information is recorded.

# **Accounting (associations)**

An Association is a combination of a Cluster, a User, and an Account.

- An accounting database may be used by multiple Clusters.
- Account is a slurm entity.
- User is a Linux user.

Use –account srun option.

With associations, a user may have different privileges on different clusters.

A user may also be able to use different accounts, with different privileges.

Multiple users may launch jobs on a linux account.

# Sacctmgr

## **Account Options**

**Clusters** to which the Account has access

Name, Description and Organization.

**Parent** is the name of an account for which this account is a child.

## **User Options**

**Account(s)** to which the user belongs.

AdminLevel is accounting privileges (for sacctmgr). None, Operator, Admin

Cluster limits clusters on which accounts user can be added to.

**DefaultAccount** is the account for the user if an account is not specified on srun

**Partition** is the a partition an association applies to.

# **Accounting Limits Enforcement**

If a user has a limit set SLURM will read in those, if not we will refer to the account associated with the job. If the account doesn't have the limit set we will refer to the cluster's limits. If the cluster doesn't have the limit set no limit will be enforced.

Some (but not all limits are)

**Fairshare**= Integer value used for determining priority. Essentially this is the amount of claim this association and it's children have to the above system. Can also be the string "parent", this means that the parent association is used for fairshare.

**GrpCPUMins=** A hard limit of cpu minutes to be used by jobs running from this association and its children. If this limit is reached all jobs running in this group will be killed, and no new jobs will be allowed to run. (GrpCPUs, GrpJobs, GrpNodes, GrpSubmitJobs, GrpWall)

MaxCPUMinsPerJob= A limit of cpu minutes to be used by jobs running from this association. If this limit is reached the job will be killed will be allowed to run. (MaxCPUsPerJob, MaxJobs, MaxNodesPerJob, MaxSubmitJobs, MaxWallDurationPerJob)

**QOS** (quality of service) comma separated list of QOS's this association is able to run.

**Important Note:** To activate the accounting limitations and QOSyou need to add the following parameter in slurm.conf, distribute the slurm.conf on all nodes and restart the deamons: AccountingStorageEnforce=limits, qos

# **Partitions and QOS**

- Partitions and QOS are used in SLURM to group nodes and jobs characteristics
- The use of Partitions and QOS (Quality of Services) entities in SLURM is orthogonal:
  - Partitions for grouping resources characteristics
  - QOS for grouping limitations and priorities

**Partition 1:** 32 cores and high\_memory

Partition 2: 32 cores and low\_memory

Partition 3: 64 cores

QOS 1:
-High priority
-Higher limits

**QOS 2:** 

-Low Priority

-Lower limits

# **Partitions and QOS Configuration**

## **Partitions Configuration:**

In slurm.conf file

#### **# Partition Definitions**

PartitionName=all Nodes=trek[0-3] Shared=NO Default=YES
PartitionName=P2 Nodes=trek[0-3] Shared=NO Priority=2 PreemptMode=CANCEL
PartitionName=P3 Nodes=trek[0-3] Shared=Exclusive Priority=3 PreemptMode=REQUEUE

## **QOS Configuration:**

In Database

>sacctmgr add qos name=lowprio priority=10 PreemptMode=Cancel GrpCPUs=10 MaxWall=60 MaxJobs=20					
>sacctmgr add qos name=hiprio priority=100 Preempt=lowprio GrpCPUs=40 MaxWall=120 MaxJobs=50					
>sacctmgr list qos					
Name Priority Preempt PreemptMode GrpCPUs MaxJobs MaxWall					

Name	Priority	reempt PreemptMode	e GrpCPU	s MaxJobs	Maxwaii 	
lowprio hiprio		cancel lowprio	10 40	20 50	60 120	

# More on QOS

Used to provide detailed limitations and prioritiees on jobs

Every user/account will have multiple allowed QOS upon which he may send jobs with the –qos parameter but only one default QOS in case he doesn't precise a –qos parameter in his submission

**Important Note:** To view the particular parameters of QOS provided by the admins users can use the "sacctmgr show associations" command

## Fairsharing in SLURM

- User and Group accounts created in the database
- Inheritance between Groups and Users for all the different characteristics (Fairshare factors, Max number of Jobs, Max number of CPUs, etc)
- Job Priorities based on the CPU\*Time utilization of each user
- Various factors can take part in the formula through the MultiFactor plugin:

```
Job_priority =

(PriorityWeightAge) * (age_factor) +

(PriorityWeightFairshare) * (fair-share_factor) +

(PriorityWeightJobSize) * (job_size_factor) +

(PriorityWeightPartition) * (partition_factor)
```

**Important Note:** To activate fairsharing in SLURM you need to add the Priority/multifactor parameter in slurm.conf along with the different parameters for the particular factors that are needed for the site

# **Usage Guide - Accounting**

# sacct displays accounting information for jobs and steps

Some basic parameters for **sacct** command:

- **-b** Displays a brief listing (jobid, status, exitcode)
- -I a long listing of jobs characteristics
- --format <param1,param2,> to select the actual fields to be shown

Example: >sacct -format=jobid, elapsed, ncpus, ntasks, state # sacctformat=jobid, elapsed, ncpus, ntasks, state					
Jobid	Elapsed	Ncpus Nt	tasks .	State	
3	00:01:30	2	2 1	COMPLETED	
3.0	00:01:30	2	2 1	COMPLETED	
4	00:00:00	2	? 2	COMPLETED	
4.0	00:00:01	2	? 2	COMPLETED	
5	00:01:23	2	2 1	COMPLETED	
5.0	00:01:31	2	2 1	COMPLETED	

# **Usage Guide - Reporting**

**sreport** generates reports of job usage and cluster utilization

The syntax of this command is like:

<type><REPORT><OPTIONS> where <type> can be cluster, job or user and each type has various reports and options

Example1: sreport job sizesbyaccount

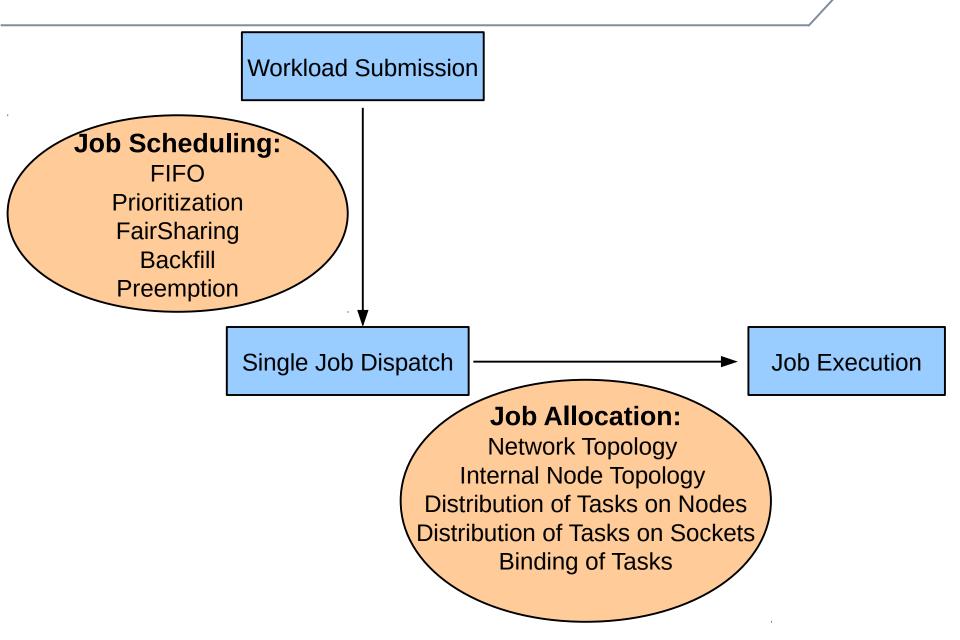
Example2: sreport cluster AccountUtilizationByUser

Example3: sreport user topusage account=gohn

Example:

>sreport cluster utilization

# **SLURM** scheduling / allocation procedures



# **SLURM Scheduling**

- SLURM supports various scheduling policies and optimization techniques (non-exhaustive list) :
  - Backfill
  - Preemption
  - Fairsharing
  - Topology aware placement
- Advantage: Techniques can be supported simultaneously
- Scheduler scalable upto 130.000 cores and 90.000 jobs per hour (confirmed upon Tera 100)
- Enhancements for High Throughput Computing
- Study experimenting with scalability and network topology-aware placement has been published in 2012:

http://link.springer.com/chapter/10.1007/978-3-642-35867-8\_8#

# **Scheduling Policies**

## **Scheduler Plugin Type**

Sched/builtin Default FIFO

Sched/hold variation on builtin; new jobs are held if

/etc/slurm.hold file exists.

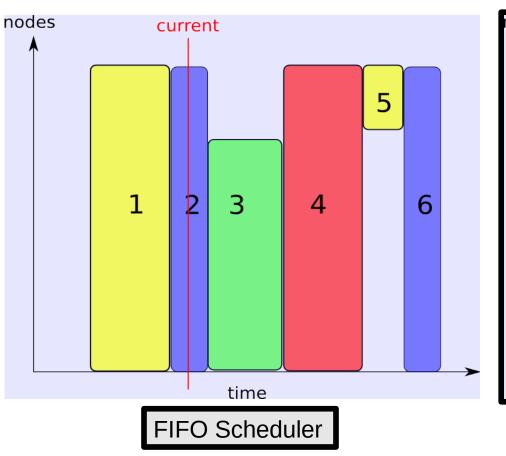
**Sched/backfill** schedule lower priority jobs as long as they don't delay a waiting higher priority job.

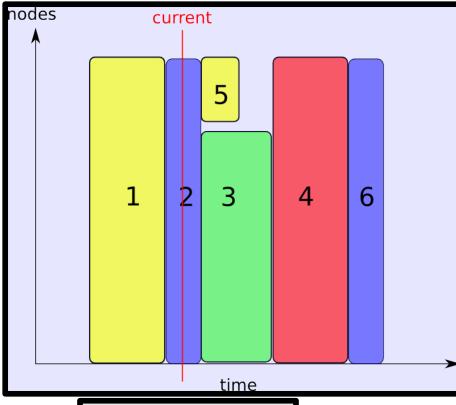
- Increases utilization of the cluster.
- Requires declaration of max execution time of lower priority jobs.
  - --time on 'srun',
  - DefaultTime or MaxTime on Partition
  - MaxWall from accounting association

#slurm.conf file SchedulerType=sched/backfill SchedulerParameters=defer,bf\_interval=60 FastSchedule=1

# Scheduling - Backfill

# Holes can be filled if previous jobs order is not changed





Backfill Scheduler

# **Scheduling Configuration Tips - Backfill**

Important parameter for **backfill** to take effect is the **Walltime** of the job (Max time allowed for the job to be completed).

- Through command line option (--time=<Minutes>)
- Partitions or QOS can be declared with Walltime parameter and jobs submitted to these partitions inherit automatically those parameters.

```
Configuration of scheduler backfill in slurm.conf

Scheduler Parameters= bf_interval=#, bf_max_job_user=#,

bf_resolution=#,bf_window=#,max_job_bf=#
```

# **Preemption Policies**

### **Preempt Types**

### None

**Partition\_prio** priority on partition definition.

**Qos** quality of service defined in accounting database.

### **Preempt Modes**

### Off

**Cancel** preempted job is cancelled.

**Checkpoint** preempted job is checkpointed if possible, or cancelled.

**Gang** enables time slicing of jobs on the same resource.

**Requeue** job is requeued as restarted at the beginning (only for sbatch).

**Suspend** job is suspended until the higher priority job ends (requires Gang).

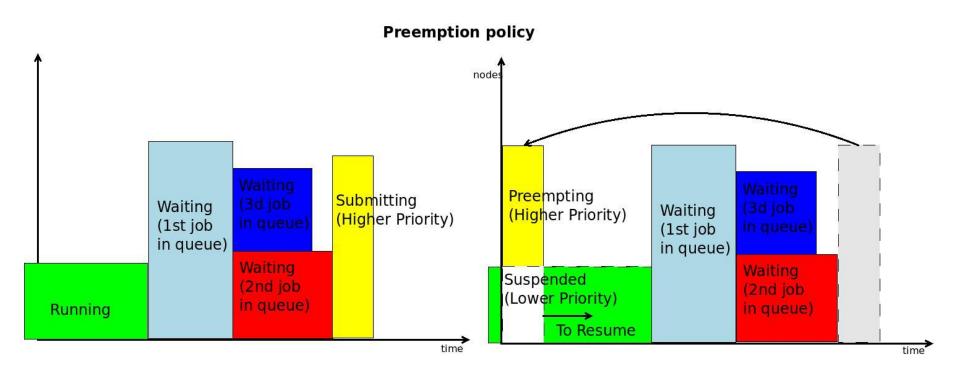
#slurm.conf file

PreemptMode=REQUEUE

PreemptType=preempt/gos

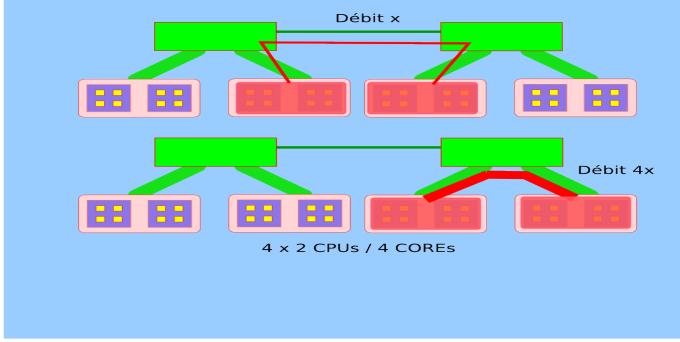
# **Scheduling - Preemption**

Preemption policy allows higher priority jobs to execute without waiting upon the cluster resources by taking the place of the lower priority jobs



# **Network Topology Aware Placement**

- topology/tree SLURM Topology aware plugin. Best-Fit selection of resources
- In fat-tree hierarchical topology: Bisection Bandwidth Constraints need to be taken into account



#slurm.conf file TopologyPlugin=topology/tree

# **Configuration (topology.conf)**

**topology.conf file** needs to exist on all computing nodes for network topology architecture description

```
# topology.conf file
SwitchName=Top
Switches=TS1,TS2,TS3,TS4,TS5,TS6,...

SwitchName=TS1 nodes=curie[1-18]
SwitchName=TS2 nodes=curie[19-37]
SwitchName=TS3 nodes=curie[38-56]
SwitchName=TS4 nodes=curie[57-75]
....
```

# **Network Topology Aware Placement**

In the slurm.conf the topology/tree plugin may be activated by the admins to allow job placement according to network topology constraints

In the submission commands the users may use the
--switches=<count>[@<max-time>] parameter to indicate
how many switches their job would be ideal to execute upon:
When a tree topology is used, this defines the maximum
count of switches desired for the job allocation and optionally
the maximum time to wait for that number of switches.

# Generic Resources (Allocation of GPUs, MIC, etc)

Generic Resources (GRES) are resources associated with a specific node that can be allocated to jobs and steps. The most obvious example of GRES use would be GPUs. GRES are identified by a specific name and use an optional plugin to provide device-specific support.

SLURM supports no generic resourses in the default configuration. One must explicitly specify which resources are to be managed in the **slurm.conf** configuration file. The configuration parameters of interest are:

- GresTypes a comma delimited list of generic resources to be managed (e.g. GresTypes=gpu,nic). This name may be that of an optional plugin providing additional control over the resources.
- **Gres** the specific generic resource and their count associated with each node (e.g. NodeName=linux[0-999] Gres=gpu:8,nic:2) specified on all nodes and SLURM will track the assignment of each specific resource on each node. Otherwise SLURM will only track a count of allocated resources rather than the state of each individual device file.

## **Generic Resources (Allocation of GPUs, MIC, etc)**

**For configuration** the new file **gres.conf** needs to exist on each compute node with gres resources

# Configure support for our four GPUs

Name=gpu File=/dev/nvidia0 CPUs=0,1

Name=gpu File=/dev/nvidia1 CPUs=0,1

Name=gpu File=/dev/nvidia2 CPUs=2,3

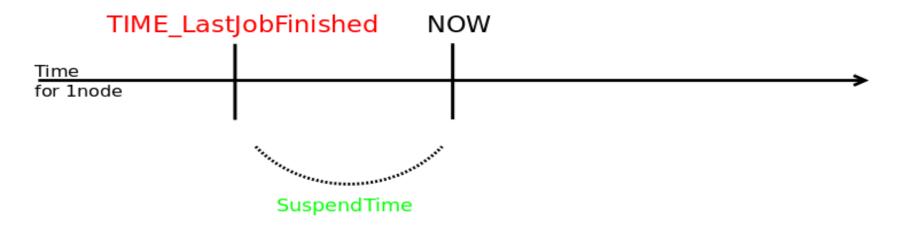
Name=gpu File=/dev/nvidia3 CPUs=2,3

**For job execution** the –gres option has to be used for to salloc, sbatch, and srun.

--gres=<list>Specifies a comma delimited list of generic consumableresources. The format of each entry on the list is "name[:count]".

# **Energy reduction techniques**

- Parameters for energy reduction techniques
- Automatic node shut-down or other actions in case of resources unutilization during particular time.



### Algorithm for SLURM Energy Reduction Techniques

Nodes Sleep Actions
if SuspendTime > A\_PreDefined\_Idle\_TIME
exec SuspendProgram upon SuspendRate nodes per minute

Nodes WakeUp Actions
if SleepingNode\_isNeeded then
exec ResumeProgram upon ResumeRate nodes per minute

# **Energy reduction techniques Configuration**

**SuspendTime:** Idle time to activate energy reduction techniques. A negative number disables power saving mode. The default value is -1 (disabled).

**SuspendRate:** # nodes added per minute. A value of zero results in no limits being imposed. The default value is 60. Use this to prevent rapid drops in power consumption.

**ResumeRate:** # nodes removed per minute. A value of zero results in no limits being imposed. The default value is 300. Use this to prevent rapid increases in power consumption.

**SuspendProgram:** Program to be executed to place nodes into power saving mode. The program executes as SlurmUser (as configured in slurm.conf). The argument to the program will be the names of nodes to be placed into power savings mode (using Slurm's hostlist expression format).

**ResumeProgram:** This program may use the scontrol show node command to insure that a node has booted and the slurmd daemon started.

SuspendTimeout, ResumeTimeout, SuspendExcNodes,SuspendExcParts, BatchStartTimeout

# Controlling jobs' time limits

Srun parameter -t, --time=<time>
Slurm.conf parameter killwait

### -t --time=<time>

Set a limit on the total run time of the job or job step. If the requested time limit for a job exceeds the partition's time limit, the job will be left in a PENDING state (possibly indefinitely). When the time limit is reached, each task in each job step is sent SIGTERM followed by SIGKILL.

#### killwait

The interval, in seconds, given to a job's processes between the SIGTERM and SIGKILL signals upon reaching its time limit. If the job fails to terminate gracefully in the interval specified, it will be forcibly terminated. default= 30sec

# **Controlling jobs' memory limits**

Srun parameters –mem=<MB>,--mem-per-cpu=<MB>
Slurm.conf paramaters DefMemPerCPU, DefMemPerNode, MaxMemPerCPU
MaxMemPerCPU

#### -mem=<MB>

Specify the real memory required per node in MB.Enforcement of memory limits currently relies upon enabling of accounting, which samples memory use on a periodic basis (data need not be stored, just collected). In both cases memory use is based upon the job's Resident Set Size (RSS). A task may exceed the memory limit until the next periodic accounting sample.

#### MaxMemPerNode

Maximum real memory size available per allocated node in MegaBytes. Used to avoid over-subscribing memory and causing paging.

# **Using different MPI libraries**

### **OpenMPI**

The system administrator must specify the range of ports to be reserved in the slurm.conf file using the MpiParams parameter. For example:

MpiParams=ports=12000-12999

Launch tasks using the srun command plus the option --resv-ports.

Alternately define the environment variable SLURM\_RESV\_PORT srun –resv-ports -n <num\_procs> a.out

If OpenMPI is configured with --with-pmi either pmi or pmi2 the OMPI jobs can be launched directly using the srun command. This is the preferred way. If the pmi2 support is enabled then the command line options '--mpi=pmi2' has to be specified on the srun command line.

srun --mpi=pmi2 -n <num\_procs> a.out

### Intel-MPI

Set the *I\_MPI\_PMI\_LIBRARY* environment variable to point to the SLURM Process Management Interface (PMI) library:

export I\_MPI\_PMI\_LIBRARY=/path/to/slurm/pmi/library/libpmi.so

Use the *srun* command to launch the MPI job:

srun -n <num procs> a.out

## **Licenses Management**

For local cluster configuration use parameter for slurm.conf Licenses

Specification of licenses (or other resources available on all nodes of the cluster) which can be allocated to jobs. License names can optionally be followed by a colon and count with a default count of one. Multiple license names should be comma separated (e.g. "Licenses=foo:4,bar").

For global cluster configuration use sacctmgr add/modify/delete resource Example: sacctmgr add resource name=flexlis cluster=nazgul count=100 percentallowed=50 server=flex\_host servertype=flexlm type=license Check with scontrol show license

For execution use srun parameter -L, --licenses=<license>

Specification of licenses (or other resources available on all nodes of the cluster) which must be allocated to this job. License names can be followed by a colon and count (the default count is one). (e.g. "--licenses=foo:4,bar").

- Introduction
- SLURM scalable and flexible RJMS
- Part 1: Basics
  - Overview, Architecture, Configuration files, Partitions, Plugins, Reservations, Reconfiguration
- Part 2: Advanced Configuration
  - Accounting, Scheduling, Allocation, Network Topology Placement, Generic Resources, Licenses Management, Energy Reduction Techniques
- Part 3: Experts Configuration
  - CPU Management, Isolation with cgroups, Power Management, Simulation and evaluation
- Upcoming Features



## **SLURM Resource Management**

- Fine Resource Management through submission command parameters
  - Adapted for MultiCore/MultiThread Architectures
  - Able to treat all kind of consumable resources (nodes, CPU hierarchies, memory, GPUs, etc)
- Based upon plpa/hwloc or Cpusets for CPUs confinement and optional tasks binding upon particular CPUs

## **CPU Management Documentation**

#### **HTML documents**

 CPU Management User and Administrator Guide (http://www.schedmd.com/slurmdocs/cpu\_management.html)

#### man pages

- slurm.conf
- · srun
- · salloc
- · sbatch
- sacctmgr (for accounting limits)

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## **Allocation Policies**

## **Select Plugin Types**

**Linear** entire nodes are allocated, regardless of the number of tasks (cpus) required.

Cons\_res cpus and memory as a consumable resource. Individual resources on a node may be allocated (not shared) to different jobs. Options to treat CPUs, Cores, Sockets, and memory as individual resources that can be independently allocated. Useful for nodes with several sockets and several cores per socket.

#slurm.conf file SelectType=select/cons\_res SelectTypeParameters=CR Core Memory,CR CORE DEFAULT DIST BLOCK

## **Task Assignment Policies**

## Task Plugin controls assignment (binding) of tasks to CPUs

**None** All tasks on a node can use all cpus on the node.

**Cgroup** cgroup subsystem is used to contain CPUs allocated to a job. Portable Hardware Locality (hwloc) library used to bind tasks to CPUs.

**Affinity** Bind tasks with one of the following <u>methods</u> **Cpusets**use cpuset subsystem to contain cpus assigned to tasks. **Sched** use sched\_setaffinity to bind tasks to cpus.

In addition to the method, a <u>binding unit</u> may also be specified. It can be one of **Sockets, Cores, Threads, None**Both the method and unit are specified on the TaskPluginParam statement.

#slurm.conf file TaskPlugin=task/affinity TaskPluginParam=Cpusets,Cores

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## **CPU Management Terminology**

#### **Allocation**

Assignment of a specific set of CPU resources (nodes, sockets, cores and/or threads) to a specific job or step

#### Distribution

- 1. Assignment of a specific task to a specific node, or
- 2. Assignment of a specific task to a specific set of CPUs within a node (used for optional Task-to-CPU binding)

## **Binding**

Confinement/locking of a specific set of tasks to a specific set of CPUs within a node

## **SLURM Nodes/CPUs allocation procedure**

SLURM uses four basic steps to manage CPU resources for a job/step:

Step 1: Selection of Nodes

**Step 2:** Allocation of CPUs from the selected Nodes

**Step 3:** Distribution of Tasks to the selected Nodes

Step 4: Optional Distribution and Binding of Tasks to

CPUs within a Node

- SLURM provides a rich set of configuration and command line options to control each step
- Many options influence more than one step
- Interactions between options can be complex and difficult to predict
- Users may be constrained by Administrator's configuration choices

## **Notable Options for Step 1: Selection of Nodes**

#### Configuration options in slurm.conf

**Nodename**: Defines a node and its characteristics. This includes the layout of sockets, cores, threads and the number of logical CPUs on the node.

**FastSchedule**: Allows administrators to define "virtual" nodes with different layout of sockets, cores and threads and logical CPUs than the physical nodes in the cluster.

**PartitionName**: Defines a partition and its characteristics. This includes the set of nodes in the partition.

#### Command line options on **srun/salloc/sbatch** commands

- **--partition, --nodelist**: Specifies the set of nodes from which the selection is made
- **-N, --nodes**: Specifies the minimum/maximum number of nodes to be selected
- -B, --sockets-per-node, --cores-per-socket, --threads-per-core: Limits node selection to nodes with the specified characteristics

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#### **Notable Options for Step 2: Allocation of CPUs from Selected Nodes**

#### Configuration options in slurm.conf:

#### SelectType:

SelectType=select/linear: Restricts allocation to whole nodes SelectType=select/cons\_res: Allows allocation of individual sockets, cores or threads as consumable resources

**SelectTypeParameters**: For select/cons\_res, specifies the consumable resource—type and default allocation method within nodes

#### Command line options on **srun/salloc/sbatch**:

- -n, --ntasks: Specifies the number of tasks. This may affect the number of CPUs allocated to the job/step
- -c, --cpus-per-task: Specifies the number of CPUs per task. This may affect the number of CPUs allocated to the job/step

## Configuration options in slurm.conf:

**MaxTasksPerNode**: Specifies maximum number of tasks per node

## Command Line options on srun/salloc/sbatch:

**-m, --distribution**: Controls the order in which tasks are distributed to nodes.

## Configuration options in slurm.conf:

## TaskPlugin:

**TaskPlugin=task/none**: Disables this step.

**TaskPlugin=task/affinity**: Enables task binding using the task affinity plugin.

**TaskPlugin=task/cgroup**: Enables task binding using the new task cgroup plugin.

**TaskPluginParam**: For task/affinity, specifies the binding unit (sockets, cores or threads) and binding method (sched\_setaffinity or cpusets)

### Command Line options on srun/salloc/sbatch:

--cpu\_bind: Controls many aspects of task affinity

-m, --distribution: Controls the order in which tasks are distributed to allocated CPUs on a node for binding

SLURM uses two default methods for allocating and distributing individual CPUs from a set of resources

- <u>block method</u>: Consume all eligible CPUs consecutively from a single resource before using the next resource in the set
- <u>cyclic method</u>: Consume eligible CPUs from each resource in the set consecutively in a round-robin fashion

The following slides illustrate the default method used by SLURM for each step.

## Different ways of selecting resources in SLURM:

- Cyclic method (Balance between nodes / Round Robin )
- Block method (Minimization of fragmentation )

•	Cyclic
	[bench@wardlaw0 ~]\$ srun -n10 -N2 -exclusive /bin/hostname
	wardlaw67
	wardlaw66

```
• Block

[bench@wardlaw0 ~]$ srun -n10 -N2 /bin/hostname

wardlaw67

wardlaw67
```

#### **Step 2a: Allocation of CPUs from a set of Nodes**

The default allocation method for this case is **block**Example: A partition contains 3 nodes. Each node has 8 available CPUs.

Node	n0	n1	n2
Number of allocated CPUs	8	6	1

Users can override this default using the appropriate command line options, e.g. --nodes, --ntasks-per-node

#### Step 2b: Allocation of CPUs from a set of Sockets within a Node

The default allocation method for this case is **cyclic** Example: A node contains 2 sockets. Each socket has 4 available CPUs.

srun --ntasks=6 ...

Socket#	0	1
Number of allocated CPUs	3	3

Administrators can change the default allocation method for this case to **block** with SelectTypeParameters=CR\_CORE\_DEFAULT\_DIST\_BLOCK

Users can override the default using the command line option --distribution

#### **Step 3: Distribution of Tasks to Nodes**

The default distribution method for this case is **block** Example: A partition has 3 nodes. Each node has 8 available cpus.

Node	n0	n1	n2
Number of allocated CPUs	8	6	2
Distribution of Tasks, by Task#	0 - 3	4 - 6	7

Users can override this default using the command line option -- distribution. The option supports three alternate methods for distributing tasks to nodes: **cyclic**, **plane** and **arbitrary**.

# <u>Step 4: Distribution of Tasks to Allocated CPUs within a Node for Task-to-CPU binding</u>

The default distribution method for this case is **cyclic** Example: A node has 2 sockets. Each socket has 4 available CPUs (cores).

Node	n0											
Socket#		(	)		1							
CPU#	0	1	2	3	4	5	6	7				
Bound Task#	0	2	4	6	1	3	5	7				

Users can override this default and specify **block** distribution of tasks to CPUs using the command line option --distribution

## **CPU Resource Sharing**

- CPU Resource Sharing between Jobs
- Shared keyword on partition definition in slurm.conf
- · --share, --exclusive options on srun/salloc/sbatch command line

#### Example:

```
Nodename=n0 CPUs=8
PartitionName=p1 Nodes=n0 Shared=YES
```

```
srun --partition=p1 --ntasks=8 --share ...
srun --partition=p1 --ntasks=8 --share ...
```

## **CPU Resource Sharing**

- CPU Resource Sharing between Tasks of the same Job
- --overcommit option on srun/salloc/sbatch command line

#### Example:

```
NodeName=n0 CPUs=8
PartitionName=p1 Nodes=n0

srun --partition=p1 --ntasks=12 --overcommit ...
```

#### Note:

It is important to understand that the Linux scheduler is not aware of CPU allocations by SLURM. Unless Task-to-CPU binding is used, Linux may run any task on any CPU on the node to which the task was distributed. In this way, CPUs may be shared between tasks and jobs even in the absence of shared CPU allocation by SLURM.

CPU management by SLURM users is subject to limits imposed by SLURM Accounting. Accounting limits may be applied on CPU usage at the level of users, associations and clusters. CPU-related accounting limits include the following:

**MaxNodes:** Maximum number of nodes that can be selected for each job.

**MaxCPUs:** Maximum number of CPUs that can be allocated to each job.

**GrpNodes:** Maximum number of nodes that can be selected for all running jobs combined in this association

**GrpCPUs:** Maximum number of CPUs that can be allocated to all running jobs combined in this association

For more details, see the sacctmgr man page.

#### <u>Information about Node Selection and CPU allocation (Steps 1 & 2)</u>

```
[sulu] (slurm) mnp> srun --nodes=2 --ntasks=3 --cpus-per-task=2 sleep 60
[2] 22841
[sulu] (slurm) mnp> squeue
                          USER ST
 JOBID PARTITION
                   NAME
                                       TIME NODES NODELIST(REASON)
   309 allnodes
                  sleep
                          slurm R
                                       0:02
                                                2 n[6-7]
[sulu] (slurm) mnp> scontrol --details show job 309
  NumNodes=2 NumCPUs=6 CPUs/Task=2 RegS:C:T=*:*:*
    Nodes=n6 CPU IDs=4-7 Mem=0
    Nodes=n7 CPU IDs=6-7 Mem=0
```

#### Note:

The CPU\_IDs reported by scontrol are SLURM abstract CPU numbers, not physical CPU numbers known to Linux.

## Getting Information About CPU Usage (cont'd)

311.0

311.0

#### <u>Information about Distribution of Tasks to Nodes (Step 3)</u>

```
[sulu] (slurm) mnp> srun --nodes=3 --ntasks=5 sleep 60 &
[sulu] (slurm) mnp> squeue
 JOBID PARTITION
                    NAME
                            USER ST
                                          TIME NODES NODELIST(REASON)
   311 allnodes
                   sleep
                            slurm
                                           0:03
                                                    3 n[6-7,13]
                                 R
[sulu] (slurm) mnp> sstat --allsteps --jobs 311 --pidformat
                         Nodelist
                                                  Pids
       JobID
311.0
                              n13
                                             7994, 7995
```

3838,3839

7199

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n6

n7

## Getting Information About CPU Usage (cont'd)

#### <u>Information about Task-to-CPU binding (Step 4)</u>

```
[sulu] (slurm) mnp> srun --partition=bones-scotty --nodes=2 --ntasks=4
--cpu_bind=cores,verbose --label cat /proc/self/status | grep
Cpus_allowed_list

0: cpu_bind=MASK - scotty, task  0  0  [4070]: mask 0x8 set
3: cpu_bind=MASK - bones, task  3  0  [23808]: mask 0x80 set
1: cpu_bind=MASK - scotty, task  1  1  [4071]: mask 0x20 set
2: cpu_bind=MASK - scotty, task  2  2  [4072]: mask 0x80 set

3: Cpus_allowed_list:  7
0: Cpus_allowed_list:  3
1: Cpus_allowed_list:  5
2: Cpus_allowed_list:  7
```

#### **Example 2: Consumable Resources with balanced allocation across Nodes**

A job requires 9 CPUs (3 tasks and 3 CPUs per task). Allocate 3 CPUs from each of the 3 nodes in the default partition.

#### In slurm.conf:

SelectType=select/cons\_res
SelectTypeParameters=CR\_Core

#### <u>srun command line</u>:

srun --ntasks=3 --cpus-per-task=3 --ntasks-per-node=1 ...

To satisfy --ntasks-per-node=1, SLURM must allocate 3 CPUs on each of the 3 nodes in the default partition. This overrides the default method for allocating CPUs from a set of nodes (block allocation).

# Example 3: Consumable Resources with minimization of resource fragmentation

A job requires 12 CPUs. Allocate CPUs using the minimum number of nodes (2 nodes) and the minimum number of sockets (3 sockets) required for the job in order to minimize fragmentation of allocated/unallocated CPUs in the cluster.

#### In slurm.conf:

SelectType=select/cons\_res
SelectTypeParameters=CR\_Core, CR\_CORE\_DEFAULT\_DIST\_BLOCK

#### srun command line:

srun --ntasks=12

The default allocation method across a selection of nodes is block. This minimizes the number of nodes used for the job. The configuration option CR\_CORE\_DEFAULT\_DIST\_BLOCK sets the default allocation method across sockets within a node to block. This minimizes the number of sockets used for the job within a node.

## Example 4: Consumable resources with Task-to-CPU binding and non-default allocation and distribution methods

A job requires 18 CPUs. Allocate 6 CPUs on each node using block allocation within nodes. Use cyclic distribution of tasks to nodes and block distribution of tasks to CPUs for binding.

#### In slurm.conf:

SelectType=select/cons\_res SelectTypeParameters=CR\_Core TaskPlugin=task/affinity TaskPluginParam=sched

#### srun command line:

```
srun --ntasks=18 --ntasks-per-node=6 --distribution=cyclic:block
--cpu_bind=cores ...
```

(continued)

- -ntasks-per-node=6 overrides the default allocation method across nodes (block) and causes SLURM to allocate 6 CPUs on each of the 3 nodes in the default partition.
- --distribution cyclic:xxxxx overrides the default method for distributing tasks to nodes (block).
- task/affinity plus --cpu\_bind=cores causes SLURM to bind each task to a single allocated core.
- --distribution xxxxx:block overrides the default allocation method within nodes (cyclic) and the default method for distributing tasks to CPUs for binding (cyclic).

The following table depicts a possible pattern of allocation, distribution and binding for this job.

(continued)

srun --ntasks=18 --ntasks-per-node=6 --distribution=cyclic:block
--cpu\_bind=cores ...

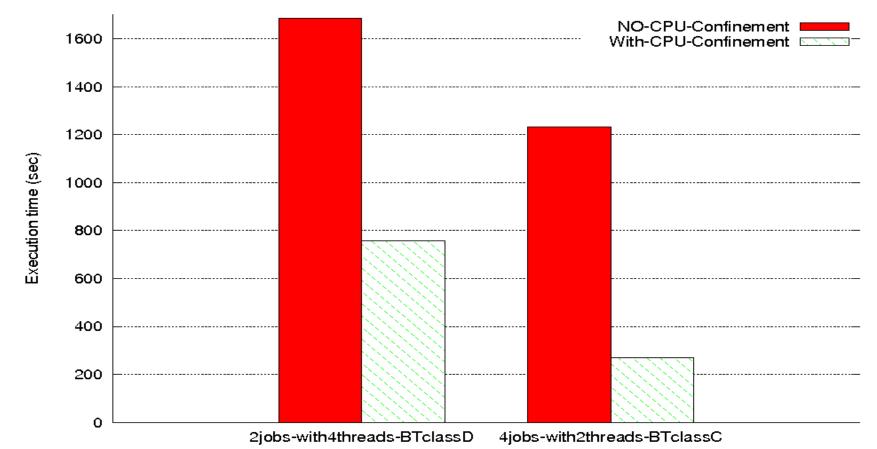
Node					n	0				n1							n2								
Socket#			C	)			1	L			(	)		1				0				1			
Numb Allocate		/1				2			4 2				4					2							
Allocated CPU#'s		0 - 5							0 - 5							0 - 5									
Number	of Tasks	6						6						6											
	ution of Nodes, ask#		0	, 3,	6, 9	9, 1	2, 1	5		1, 4, 7, 10, 13, 1					<u>.</u> 6			2,	, 5,	8, 1	.1, 1	.4, 1	L7		
Task-to-	CPU#	0	1	2	3	4	5	6	7	0	1	2	3	4	5	6	7	0	1	2	3	4	5	6	7
Binding	CPU Binding Task#		3	6	9	12	15	-	-	1	4	7	10	13	16	-	-	2	5	8	11	14	17	-	-

#### Issues with no Task Confinement upon the allocated resources

- SMP system without some means of CPU placement, any task can run on any CPU.
  - This may cause CPU idleness while other CPUs are shared and system time spent on migrating tasks between processors
- NUMA system, any memory page can be allocated on any node.
   This can cause both poor cache locality and poor memory access times.

# Execution of NAS benchmarks (MPI-OpenMP) upon 8 nodes SMP cluster (2sockets-4cores/socket) with and without CPU confinement

Simultaneous execution of hybrid OpenMP-MPI jobs (NAS benchmark BT-MZ16 processes used) upon the same 8 nodes (2sockets-4cores/socket) with and without CPU confinement



# CPU instant utilization during execution of 2 BT-MZ jobs with 16 tasks and 4 threads per task without CPU confinement

```
top - 13:32:36 up 1:21, 1 user, load average: 7.05, 5.18, 6.79
Tasks: 256 total, 9 running, 247 sleeping, 0 stopped, 0 zombie
Cpu0 (100.0%u)s, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
Cpu1 :100.0%us, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
Cpu2 :100.0%us, 0.0%sy, 0.0%ni, 0.0%ia, 0.0%wa, 0.070mi
                                                                                          CPUs are shared
Cpu3 \100-0%vs, 0.0%sy, 0.0%ni, 0.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
                                                                                             between jobs
Cpu4 \( \delta \) \( \text{0.0\% a}, \ 0.0\% sy, \ 0.0\% ni, 100.0\% id, \ 0.0\% wa, \ 0.0\% hi, \ 0.0\% si, \ 0.0\% st
Cpu5/: 0.0%us\0.0%sy, 0.0%ni,100.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
Cpu6: 0.0%us, 0.0%sy, 0.0%si, 100.0%id, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
Cpu7 \: 0.0%us/ 0.0%sy, 0.0%ni,100.0%us, 0.0%wa, 0.0%hi, 0.0%si, 0.0%st
Mem: 18395656k total, 3346396k used, 15049260k free, 364k buffers
Swap: 1022752k total,
                        0k used, 1022752k free, 114104k cacheu
             PR NI VIRT RES SHR S %CPU MMEM TIME+ P COMMAND
 PID USER
3582 georgioy 20 0 920m 713m 3956 R 54.8 4.0 0:45.34 3 bt-mz.D.16
                                                                                              While there
3581 georgioy 20 0 921m 717m 4192 R 52,5 4.0 0:45.47 1 bt-mz .16
                                                                                              are idle CPUs
3592 georgioy 20 0 921m 713m 3924 R 51 2 4.0 0:43.02 2 st-mz.D.16
3577 georgioy 20 0 920m 717m 3940 R 50 8 4.0 0:44.81 0 bt-mz.D.16
3578 georgioy 20 0 921m 713m 3924 R 50 2 4.0 0:45.37 3 bt-mz.D.16
3594 georgioy 20 0 920m 717m 3940 R 48.5 4.0 0:43.48 0 bt-mz.D.16
3598 georgioy 20 0 921m 717m 4192 R 48.2 4.0 0:43.14 1 bt-mz.D.16
3597 georgioy 20 0 920m 713m 3956 R 43.9, 4.0 0:43.18 2 bt-mz.D.16
  1 root
          20 0 21336 1548 1280 S 0.0 0.0 9:03.60 5 init
  2 root
          20 0 0 0 S 0.0 0.0 0:00.00 5 kthreadd
```

## Advantages: cgroups support for HPC

- To guarantee that every consumed resources is consumed the way it's planned to be
  - leveraging Linux latest features in terms of process control and resource management
  - Enabling node sharing
- While enhancing the connection with Linux systems
  - Improve tasks isolation upon resources
  - Improve **efficiency** of resource management activities (e.g., process tracking, collection of accounting statistics)
  - Improve robustness (e.g. more reliable cleanup of jobs)
- And simplifying the addition of new controlled resources and features
  - prospective management of network and I/O as individual resources

## Introduction to cgroups

Control Groups (cgroups) is a **Linux kernel mechanism** (appeared in 2.6.24) to limit, isolate and monitor resource usage (CPU, memory, disk I/O, etc.) of groups of processes.

## **Features**

- •Resource Limiting (i.e. not to exceed a memory limit)
- •Prioritization (i.e. groups may have larger share of CPU)
- •Isolation (i.e. isolate GPUs for particular processes)
- •Accounting (i.e. montior resource usage for processes)
- •Control (i.e. suspending and resuming processes)

## **Cgroups Model and Concepts**

#### Model

Cgroups **similar** to Linux processes:

- Hierarchical
- Inheritance of attributes from parent to child

but **different** because:

 multiple hierarchies of cgroups may exist that are attached to one or more subsystems

#### **Concepts**

- Cgroup a group of processes with the same characteristics
- Subsystem a module that applies parameters to a group of processes (cgroup)
- Hierarchy a set of cgroups organized in a tree, plus one or more subsystems associated with that tree

## **Cgroups subsystems**

- cpuset assigns tasks to individual CPUs and memory nodes in a cgroup
- •cpu schedules CPU access to cgroups
- cpuacct reports CPU resource usage of tasks of a cgroup
- •memory set limits on memory use and reports memory usage for a cgroup
- •devices allows or denies access to devices (i.e. gpus) for tasks of a cgroup
- •freezer suspends and resumes tasks in a cgroup
- •net\_cls tags network packets in a cgroup to allow network traffic priorities
- •ns namespace subsystem
- •blkio tracks I/O ownership, allowing control of access to block I/O resources

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## **Cgroups functionality rules**

- Cgroups are represented as virtual file systems
  - Hierarchies are directories, created by mounting subsystems, using the mount command; subsystem names specified as mount options
  - Subsystem parameters are represented as files in each hierarchy with values that apply only to that cgroup
- •Interaction with cgroups take place by manipulating directories and files in the cgroup virtual file system using standard shell commands and system calls (mkdir, mount, echo, etc)
  - tasks file in each cgroup directory lists the tasks (pids) in that cgroup
  - Tasks are automatically removed from a cgroup when they terminate or are added to a different cgroup in the same hierarchy
  - Each task is present in only one cgroup in each hierarchy
- Cgroups have a mechanism for automatic removal of abandoned cgroups (release\_agent)

## **Cgroups subsystems parameters**

#### cpuset subsystem

**cpuset.cpus**: defines the set of cpus that the tasks in the cgroup are allowed to execute on

**cpuset.mems**: defines the set of memory zones that the tasks in the cgroup are allowed to use

#### memory subsystem

**memory.limit\_in\_bytes**: defines the memory limit for the tasks in the cgroup **memory.swappiness**: controls kernel reclamation of memory from the tasks in the cgroup (swap priority)

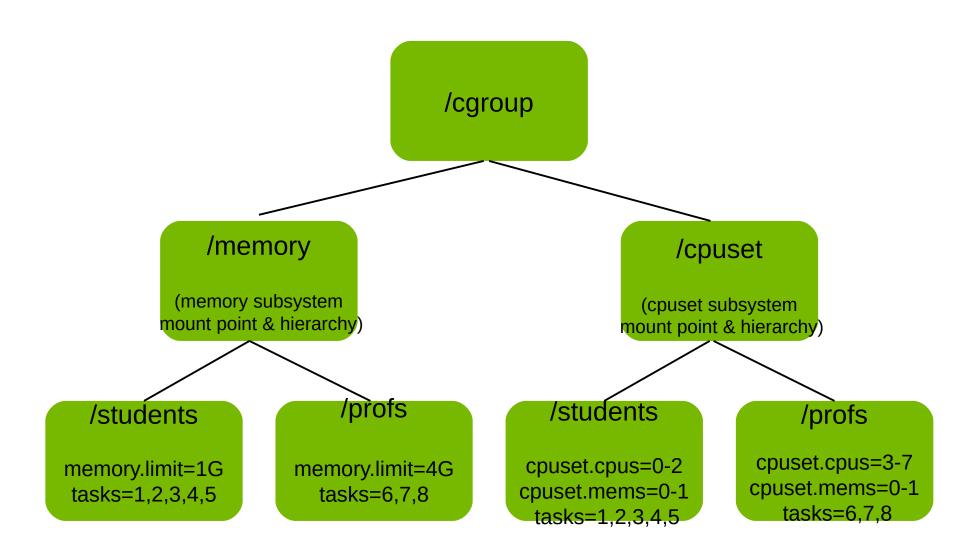
#### freezer subsystem

**freezer.state**: controls whether tasks in the cgroup are active (runnable) or suspended

#### devices subsystem

devices\_allow: specifies devices to which tasks in a cgroup have acces

#### **Cgroups functionality example**



#### **Cgroups functionality example**

```
[root@mordor:~]# mkdir /cgroup
[root@mordor:~]# mkdir /cgroup/cpuset
[root@mordor:~]# mount -t cgroup -o cpuset none /cgroup/cpuset
[root@mordor:~]# ls /cgroup/cpuset/
cpuset.cpus cpuset.mems tasks notify_on_release release_agent
[root@mordor:~]# mkdir /cgroup/cpuset/students
[root@mordor:~]# mkdir /cgroup/cpuset/profs
[root@mordor:~]# echo 0-2 > /cgroup/cpuset/students/cpuset.cpus
[root@mordor:~]# echo 0 > /cgroup/cpuset/students/cpuset.mems
[root@mordor:~]# echo $PIDS_st > /cgroup/cpuset/students/tasks
[root@mordor:~]# echo 3-7 > /cgroup/cpuset/profs/cpuset.cpus
[root@mordor:~]# echo 1 > /cgroup/cpuset/profs/cpuset.mems
[root@mordor:~]# echo $PIDS_pr > /cgroup/cpuset/profs/tasks
```

### **Process Tracking with Cgroups**

#### Track job processes using the <u>freezer</u> subsystem

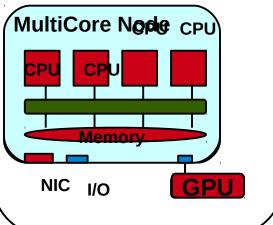
- Every spawned process is tracked
  - Automatic inheritance of parent's cgroup
  - No way to escape the container
- Every processes can be frozen
  - Using the Thawed | Frozen state of the subsystem
  - No way to avoid the freeze action

### [mat@leaf slurm]\$ srun sleep 300

```
[root@leaf ~] # cat /cgroup/freezer/uid 500/job 53/step 0/freezer.state
THAWED
[root@leaf ~]# scontrol suspend 53
[root@leaf \sim] \# ps -ef f \mid tail -n 2
root 15144 1 0 17:10? Sl 0:00 slurmstepd: [53.0]
mat 15147 15144 0 17:10 ? T 0:00 \_/bin/sleep 300
[root@leaf ~] # cat /cgroup/freezer/uid 500/job 53/step 0/freezer.state
FREEZING
[root@leaf ~]# scontrol resume 53
[root@leaf \sim] \# ps -ef f \mid tail -n 2
root 15144 1 0 17:10? Sl 0:00 slurmstepd: [53.0]
mat 15147 15144 0 17:10 ? S 0:00 \ /bin/sleep 300
[root@leaf ~] # cat /cgroup/freezer/uid 500/job 53/step 0/freezer.state
THAWED
[root@leaf ~]#
```

#### Task confinement for allocated resources

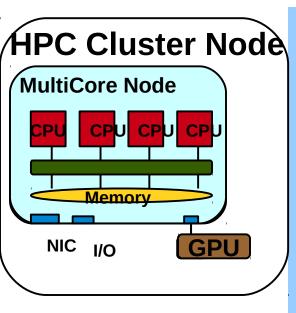
### HPC Cluster Node



#### Constrain jobs tasks to the allocated resources

- 3 independant layers of managed resources using 3 subsystems
  - Cores (<u>cpuset</u>), Memory (<u>memory</u>),
     GRES (<u>devices</u>)
- Every spawned process is tracked
  - Automatic inheritance of parent's cgroup
  - No escape, no way to use additional resources,
- Each layer has its own additional parameters
- More resources could be added in the future

#### Task confinement for cpus



#### Constrain jobs tasks to the allocated cores

- Configurable feature
  - ConstrainCores=yes no
- Use step's allocated cores with "exclusive steps"
  - Otherwise, let steps use job's allocated cores
- Basic affinity management as a configurable sub-feature
  - TaskAffinity=yes|no in cgroup.conf (rely on HWLOC)
  - Automatic block and cyclic distribution of tasks

[mat@leaf slurm]\$ salloc --exclusive srun -n1 --cpu\_bind=none sleep 3000

salloc: Granted job allocation 55

```
[root@leaf ~]# egrep "Cores | Affinity" /etc/slurm/cgroup.conf
ConstrainCores=yes
TaskAffinity=yes
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log | grep task/cgroup
[2011-09-16T17:24:59] [55.0] task/cgroup: now constraining jobs allocated
cores
[2011-09-16T17:24:59] [55.0] task/cgroup: loaded
[2011-09-16T17:24:59] [55.0] task/cgroup: job abstract cores are '0-31'
[2011-09-16T17:24:59] [55.0] task/cgroup: step abstract cores are '0-31'
[2011-09-16T17:24:59] [55.0] task/cgroup: job physical cores are '0-31'
[2011-09-16T17:24:59] [55.0] task/cgroup: step physical cores are '0-31'
[2011-09-16T17:24:59] [55.0] task/cgroup: task[0] is requesting no affinity
```

[mat@leaf slurm]\$ salloc --exclusive srun -n1 --cpu\_bind=cores sleep 3000 salloc: Granted job allocation 57

```
[root@leaf ~]# egrep "Cores | Affinity" /etc/slurm/cgroup.conf
ConstrainCores=yes
TaskAffinity=yes
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log | grep task/cgroup
[2011-09-16T17:31:17] [57.0] task/cgroup: now constraining jobs allocated cores
[2011-09-16T17:31:17] [57.0] task/cgroup: loaded
[2011-09-16T17:31:17] [57.0] task/cgroup: job abstract cores are '0-31'
[2011-09-16T17:31:17] [57.0] task/cgroup: step abstract cores are '0-31'
[2011-09-16T17:31:17] [57.0] task/cgroup: job physical cores are '0-31'
[2011-09-16T17:31:17] [57.0] task/cgroup: step physical cores are '0-31'
[2011-09-16T17:31:17] [57.0] task/cgroup: task[0] is requesting core level binding
[2011-09-16T17:31:17] [57.0] task/cgroup: task[0] using Core granularity
[2011-09-16T17:31:17] [57.0] task/cgroup: task[0] taskset '0x00000001' is set
```

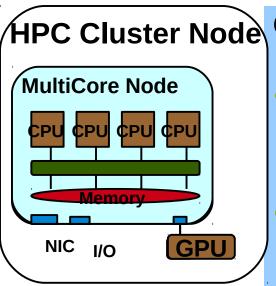
[mat@leaf slurm]\$ salloc --exclusive srun -n1 --cpu\_bind=socket sleep 3000 salloc: Granted job allocation 58

```
[root@leaf ~]# egrep "Cores | Affinity" /etc/slurm/cgroup.conf
ConstrainCores=yes
TaskAffinity=yes
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log | grep task/cgroup
[2011-09-16T17:33:31] [58.0] task/cgroup: now constraining jobs allocated cores
[2011-09-16T17:33:31] [58.0] task/cgroup: loaded
[2011-09-16T17:33:31] [58.0] task/cgroup: job abstract cores are '0-31'
[2011-09-16T17:33:31] [58.0] task/cgroup: step abstract cores are '0-31'
[2011-09-16T17:33:31] [58.0] task/cgroup: job physical cores are '0-31'
[2011-09-16T17:33:31] [58.0] task/cgroup: step physical cores are '0-31'
[2011-09-16T17:33:31] [58.0] task/cgroup: task[0] is requesting socket level binding
[2011-09-16T17:33:31] [58.0] task/cgroup: task[0] using Socket granularity
[2011-09-16T17:33:31] [58.0] task/cgroup: task[0] taskset '0x00000003' is set
```

[mat@leaf slurm]\$ salloc --exclusive srun -n2 --cpu\_bind=socket sleep 3000 salloc: Granted job allocation 60

```
[root@leaf ~]# egrep "Cores | Affinity" /etc/slurm/cgroup.conf
ConstrainCores=yes
TaskAffinity=yes
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log | grep task/cgroup[2011-09-16T17:36:18] [60.0]
task/cgroup: now constraining jobs allocated cores
[2011-09-16T17:36:18] [60.0] task/cgroup: loaded
[2011-09-16T17:36:18] [60.0] task/cgroup: job abstract cores are '0-31'
[2011-09-16T17:36:18] [60.0] task/cgroup: step abstract cores are '0-31'
[2011-09-16T17:36:18] [60.0] task/cgroup: job physical cores are '0-31'
[2011-09-16T17:36:18] [60.0] task/cgroup: step physical cores are '0-31'
[2011-09-16T17:36:18] [60.0] task/cgroup: task[0] is requesting socket level binding
[2011-09-16T17:36:18] [60.0] task/cgroup: task[1] is requesting socket level binding
[2011-09-16T17:36:18] [60.0] task/cgroup: task[1] using Core granularity
[2011-09-16T17:36:18] [60.0] task/cgroup: task[1] higher level Socket found
[2011-09-16T17:36:18] [60.0] task/cgroup: task[1] taskset '0x00000003' is set
[2011-09-16T17:36:18] [60.0] task/cgroup: task[0] using Core granularity
[2011-09-16T17:36:18] [60.0] task/cgroup: task[0] higher level Socket found
[2011-09-16T17:36:18] [60.0] task/cgroup: task[0] taskset '0x00000003' is set
```

## Task confinement for memory : **memory** subsystem



## Constrain jobs tasks to the allocated amount of memory

- Configurable feature
  - ConstrainRAMSpace=yes | no
  - ConstrainSwapSpace=yes|no
- Use step's allocated amount of memory with "exclusive steps"
  - Else, let steps use job's allocated amount
- Both RSS and swap are monitored
- Trigger OOM killer on the cgroup's tasks when reaching limits
- Tolerant mechanism
  - AllowedRAMSpace , AllowedSwapSpace percents

#### Cgroup Task plugin : memory subsystem

[mat@leaf slurm]\$ salloc --exclusive --mem-per-cpu 100 srun -n1 sleep 3000 salloc: Granted job allocation 67

```
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log |grep task/cgroup
[2011-09-16T17:55:20] [67.0] task/cgroup: now constraining jobs allocated memory
[2011-09-16T17:55:20] [67.0] task/cgroup: loaded
[2011-09-16T17:55:20] [67.0] task/cgroup: job mem.limit=3520MB memsw.limit=3840MB
[2011-09-16T17:55:20] [67.0] task/cgroup: step mem.limit=3520MB memsw.limit=3840MB
```

[mat@leaf slurm]\$ salloc --exclusive --mem-per-cpu 100 srun – exclusive -n1 sleep 3000

salloc: Granted job allocation 68

```
[root@leaf ~]# tail -f /var/log/slurmd.leaf10.log | grep task/cgroup

[2011-09-16T17:57:31] [68.0] task/cgroup: now constraining jobs allocated memory

[2011-09-16T17:57:31] [68.0] task/cgroup: loaded

[2011-09-16T17:57:31] [68.0] task/cgroup: job mem.limit=3520MB memsw.limit=3840MB

[2011-09-16T17:57:31] [68.0] task/cgroup: step mem.limit=110MB memsw.limit=120MB
```

### Cgroup **Task** plugin : **memory** subsystem OOM killer usage

[mat@leaf slurm]\$ salloc --exclusive --mem-per-cpu 100 srun -n1 sleep 3000

salloc: Granted job allocation 67

slurmd[berlin27]: Step 268.0 exceeded 1310720 KB memory limit, being killed

srun: Exceeded job memory limit

srun: Job step aborted: Waiting up to 2 seconds for job step

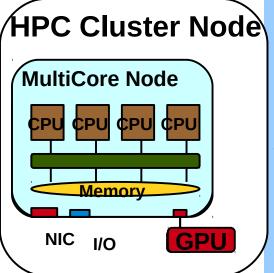
to finish.

slurmd[berlin27]: \*\*\* STEP 268.0 KILLED AT 2012-03-

31T15:50:36 WITH SIGNAL 9 \*\*\*

srun: error: berlin27: tasks 0,1: Killed

## Tasks confinement for devices: **devices** subsystem



## Constrain jobs tasks to the allocated system devices

- Based on the GRES plugin for generic resources allocation (NIC, GPUs, etc) and built upon the cgroup task plugin
  - Each task is allowed to access to a number of devices by default
  - Only the tasks that have granted allocation on the GRES devices will be allowed to have access on them.
  - Tasks with no granted allocation upon GRES devices will not be able to use them.

#### Cgroup **Task** plugin : **devices** subsystem

### **Cgroup Devices Configuration Example**

[root@mordor cgroup]# egrep "Devices" /etc/slurm/cgroup.conf ConstrainDevices=yes AllowedDevicesFile="/etc/slurm/allowed\_devices.conf"

[root@mordor cgroup]# cat /etc/slurm/allowed\_devices.conf

/dev/sda\*

/dev/null

/dev/zero

/dev/urandom

/dev/cpu/\*/\*

#### Cgroup Task plugin : devices subsystem

#### **Cgroup Devices Logic as implemented in task plugin**

- 1) Initialization phase (information collection gres.conf file, major, minor, etc)
- 2) Allow all devices that should be allowed by default (allowed\_devices.conf)
- 3) Lookup which gres devices are allocated for the job
- Write allowed gres devices to devices.allow file
- Write denied gres devices to devices.deny file
- 4) Execute 2 and 3 for job and steps tasks (different hierarchy level in cgroups)

#### Cgroups devices subsystem : Usage Example

[root@mordor cgroup]# egrep "Gres" /etc/slurm/slurm.conf GresTypes=gpu NodeName=cuzco[57,61] Gres=gpu:2 Procs=8 Sockets=2 CoresPerSocket=4

```
[root@cuzco51]# cat /etc/slurm/allowed_devices.conf /dev/sda* /dev/null
```

```
[gohn@cuzco0]$ cat gpu_test.sh
#!/bin/sh
sleep 10
echo 0 > /dev/nvidia0
echo 0 > /dev/nvidia1
```

#### Cgroups devices subsystem : Usage Example

#### [gohn@cuzco0]\$ srun -n1 -gres=gpu:1 -o output ./gpu\_test.sh

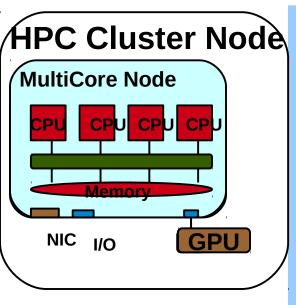
```
[root@cuzco51 ~]# tail -f /var/log/slurmd.cuzco51.log
[2011-09-20T03:10:02] [22.0] task/cgroup: manage devices jor job '22'
[2011-09-20T03:10:02] [22.0] device : /dev/nvidia0 major 195, minor 0
[2011-09-20T03:10:02] [22.0] device: /dev/nvidia1 major 195, minor 1
[2011-09-20T03:10:02] [22.0] device : /dev/sda2 major 8, minor 2
[2011-09-20T03:10:02] [22.0] device : /dev/sda1 major 8, minor 1
[2011-09-20T03:10:02] [22.0] device : /dev/sda major 8, minor 0
[2011-09-20T03:10:02] [22.0] device : /dev/null major 1, minor 3
[2011-09-20T03:10:02] [22.0] Default access allowed to device b 8:2 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.allow' set to 'b 8:2 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
[2011-09-20T03:10:02] [22.0] Default access allowed to device b 8:1 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.allow' set to 'b 8:1 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
[2011-09-20T03:10:02] [22.0] Default access allowed to device b 8:0 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.allow' set to 'b 8:0 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
[2011-09-20T03:10:02] [22.0] Default access allowed to device c 1:3 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.allow' set to 'c 1:3 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
[2011-09-20T03:10:02] [22.0] Allowing access to device c 195:0 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.allow' set to 'c 195:0 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
[2011-09-20T03:10:02] [22.0] Not allowing access to device c 195:1 rwm
[2011-09-20T03:10:02] [22.0] parameter 'devices.deny' set to 'c 195:1 rwm' for '/cgroup/devices/uid 50071/job 22/step 0'
```

#### Cgroups devices subsystem : Usage Example

```
[root@cuzco51 ~]# cat /cgroup/devices/uid_50071/job_22/step_0/tasks
4875
4879
4882
[root@cuzco51 ~]# cat /cgroup/devices/uid_50071/job_22/step_0/devices.list
b 8:2 rwm
b 8:1 rwm
b 8:0 rwm
c 1:3 rwm
c 195:0 rwm
```

### [gohn@cuzco0]\$ cat output /home/GPU/./gputest.sh: line 4: echo: write error: Invalid argument /home/GPU/./gputest.sh: line 5: /dev/nvidia1: Operation not permitted

## Monitoring Resource Usage: cpuacct and memory subsystems



### Monitoring cpu usage with cpuacct subsystem and memory usage with memory subsystem

- Implemented as a jobacct\_gather plugin for SLURM
- Collects information concerning CPU time and Memory RSS consumed for each task of the cgroup
- Values reported as a new job characteristics in the accounting database of SLURM
- Values can be used for billing purposes
- Monitor per job energy consumption (not through cgroups )

### Monitoring Resources: cpuacct -memory subsystems

[gohn@cuzco0]\$ srun -n32 ./malloc [gohn@cuzco0]\$ sacct -j 167

JobID JobName Partition MaxRSS AveRSS MaxPages AvePages MinCPU AveCPU Elapsed State Ntasks AllocCPUs ExitCode

167.0 malloc shared 61311K 57221K 239.24G 99893120K 00:03.000 00:03.000 00:01:10 COMPLETED 32 32 0.0

#### Cgroup Task plugin : devices subsystem

#### **Cgroup Devices Logic as implemented in task plugin**

- 1) Initialization phase (information collection gres.conf file, major, minor, etc)
- 2) Allow all devices that should be allowed by default (allowed\_devices.conf)
- 3) Lookup which gres devices are allocated for the job
- Write allowed gres devices to devices.allow file
- Write denied gres devices to devices.deny file
- 4) Execute 2 and 3 for job and steps tasks (different hierarchy level in cgroups)

### **Energy accounting and control**

#### Summary of the energy accounting and control features

- Power and Energy consumption monitoring per node level.
- Energy consumption accounting per step/job on SLURM DataBase
- Power profiling per step/job on the end of job
- Frequency Selection Mechanisms for user control of job energy consumption

#### Summary of the energy accounting and control features

- Power and Energy consumption monitoring per node level.
- Energy consumption accounting per step/job on SLURM DataBase
- Power profiling per step/job on the end of job
- Frequency Selection Mechanisms for user control of job energy consumption

#### How this takes place:

- Dedicated Plugins for Support of in-band collection of energy/power data (IPMI / RAPL)
- Dedicated Plugins for Support of out-of-band collection of energy/power data (RRD databases)
- Power data job profiling with HDF5 file format
- SLURM Internal power-to-energy and energy-to-power calculations

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Overhead: In-band Collection

How this \*Precision: of the measurements and internal

calculations

data (IPM

Dedicated •Scalability: Out-of band Collection

y/power

- Dedicated Plugins for Support of out-of-band collection of energy/power data (RRD databases)
- Power data job profiling with HDF5 file format
- SLURM Internal power-to-energy and energy-to-power calculations

#### In-band collection of power/energy data with IPMI

- **IPMI** is a message-based, hardware-level interface specification (may operate in-band or out-of-band)
- Communication with the Baseboard Management Controller BMC which is a specialized microcontroller embedded on the motherboard of a computer
- SLURM support is based on the FreeIPMI API:
  - http://www.gnu.org/software/freeipmi/
    - FreeIPMI includes a userspace driver that works on most motherboards without any required driver.
    - No thread interferes with application execution
- The data collected from IPMI are currently instantaneous measures in Watts
- SLURM individual polling frequency (>=1sec)
  - direct usage for power profiling
  - but internal SLURM calculations for energy reporting per job

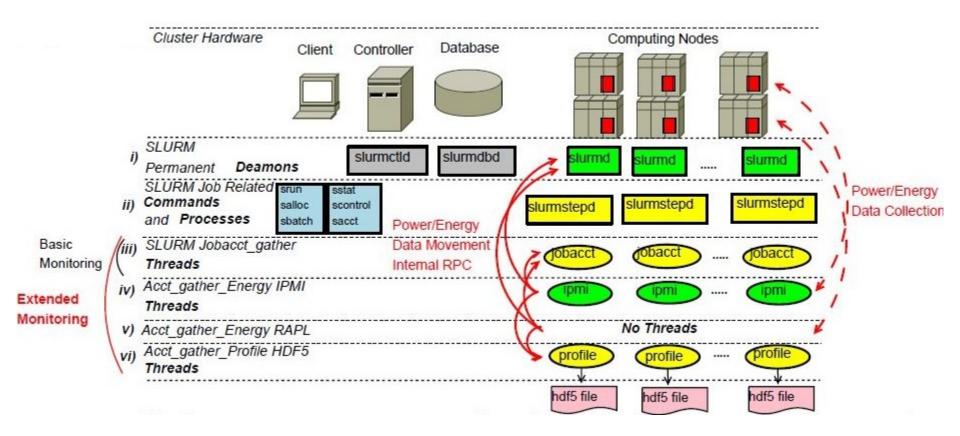
#### In-band collection of power/energy data with RAPL

- RAPL (Running Average Power Limit) are particular interfaces on Intel Sandy Bridge processors (and later models) implemented to provide a mechanism for keeping the processors in a particular user-specified power envelope.
- Interfaces can estimate current energy usage based on a software model driven by hardware performance counters, temperature and leakage models
  - Linux supports an 'MSR' driver and access to the register can be made through /dev/cpu/\*/msr with priviledged read permissions
- The data collected from RAPL is energy consumption in Joules (since the last boot of the machine)
- SLURM individual polling frequency (>=1sec)
  - direct usage for energy reporting per job
  - but internal SLURM calculations for power reporting

#### **Power Profiling**

- Job profiling to periodically capture the task's usage of various resources like CPU, Memory, Lustre, Infiniband and Power per node
- Resource Independent polling frequency configuration
- Based on hdf5 file format http://www.hdfgroup.org opensource software library
  - versatile data model that can represent very complex data objects and a wide variety of metadata
  - portable file format with no limit on the number or size of data objects stored
- Profiling per node (one hdf5 file per job on each node)
- Aggregation on one hdf5 file per job (after job termination)
- Slurm built-in tools for extraction of hdf5 profiling data

# Energy Accounting and Power Profiling Architecture



#### acct\_gather\_energy Plugin - Overview

- One of a new family of acct\_gather plugins that collect resource usage data for accounting, profiling and monitoring.
- Loaded by slurmd on each compute node.
- Called by jobacct\_gather plugin to collect energy consumption accounting data for jobs and steps.
- Called separately via RPC from the slurmctld background thread to collect energy consumption data for nodes.
- Calls acct\_gather\_profile plugin to provide energy data samples for profiling.

#### acct\_gather\_energy Plugin - Configuration

#### In slurm.conf

```
To configure plugin:
```

```
AcctGatherEnergyType=acct_gather_energy/rapl or AcctGatherEnergyType=acct_gather_energy/ipmi
```

Frequency of node energy sampling controlled by:

AcctGatherNodeFreq=<seconds>

Default value is 0, which disables node energy sampling

Collection of energy accounting data for jobs/steps requires:

JobAcctGatherType=jobacct\_gather/linux or

JobAcctGatherType=jobacct\_gather/cgroup

Frequency of job accounting sampling controlled by:

JobAcctGatherFrequency=task=<seconds>

Default value is 30 seconds

#### In acct\_gather.conf (new config file), for acct\_gather\_energy/ipmi only:

**EnergyIPMIFrequency** 

**EnergyIPMICalcAdjustment** 

**EnergyIPMIPowerSensor** 

**EnergyIPMIUsername** 

**EnergyIPMIPassword** 

#### acct\_gather\_energy Plugin - Data Reporting

- For running jobs, energy accounting data is reported by sstat.
- If accounting database is configured, energy accounting data is included in accounting records and reported by sacct and sreport.
- If acct\_gather\_profile plugin is configured, energy profiling data is reported by the method specified by the profile plugin type.
- Energy consumption data for nodes is reported by scontrol show node.
- Cumulative/total energy consumption is reported in units of joules.
- Instantaneous rate of energy consumption (power) is reported in units of watts.

#### Out-of-band collection of power/energy data

- External Sensors Plugins to allow out-of-band monitoring of cluster sensors
- Possibility to Capture energy usage and temperature of various components (nodes, switches, rack-doors, etc)
- Framework generic but initial Support for RRD databases through rrdtool API (for the collection of energy/temperature data)
  - Plugin to be used with real wattmeters or out-of-band IPMI capturing
- Power data captured used for per node power monitoring (scontrol show node) and per job energy accounting (Slurm DB)
  - direct usage for energy reporting per job
  - but internal SLURM calculations for power reporting

#### **External Sensors Plugin - Purpose**

**Plugin Name**: ext\_sensors

**Purpose**: To collect environmental-type data from external sensors or sources for the following uses:

- Job/step accounting Total energy consumption by a completed job or step (no energy data while job/step is running).
- Hardware monitoring Instantaneous and cumulative energy consumption for nodes; instantaneous temperature of nodes.
- Future work will add additional types of environmental data, such as energy and temperature data for network switches, cooling system, etc.
   Environmental data may be used for resource management.

#### ext\_sensors Plugin - Overview

- Loaded by slurmctld on management node.
- Collects energy accounting data for jobs and steps independently of the acct\_gather plugins.
  - Called by slurmctld request handler when step starts.
  - Called by slurmctld step manager when step completes.
- Since energy use by jobs/steps is measured only at completion (i.e., no sampling), does not support energy profiling or energy reporting for running jobs/steps (sstat).
- Called separately from the slurmctld background thread to sample energy consumption and temperature data for nodes.

#### ext\_sensors Plugin - Data Reporting

- If accounting database is configured, energy data is included in accounting records and reported by **sacct** and **sreport**.
- Energy consumption data for nodes is reported by scontrol show node.
- Cumulative/total energy consumption reported in joules.
- Instantaneous energy consumption rate (power) for nodes reported in watts.
- Node temperature reported in celsius.

#### ext\_sensors Plugin - Versions

- One version of ExtSensorsType plugin currently supported:
  - ext\_sensors/rrd

External sensors data is collected using RRD. RRDtool is GNU-licensed software that creates and manages a linear database used for sampling or logging. The database is populated with energy data using out-of-band IPMI collection.

- Plugin API is described in Slurm developer documentation:
  - http://slurm.schedmd.com/ext\_sensorsplugins.html

#### ext\_sensors Plugin - Configuration

In slurm.conf

To configure plugin:

ExtSensorsType=ext\_sensors/rrd

Frequency of node energy sampling controlled by:

ExtSensorsFreq=<seconds>

Default value is 0, which disables node energy sampling

Collection of energy accounting data for jobs/steps requires:

JobAcctGatherType=jobacct\_gather/linux or cgroup

In ext\_sensors.conf (new configuration file)

**JobData=energy** Specify the data types to be collected by the plugin for jobs/steps.

NodeData=[energy|temp]Specify the data types to be collected by the plugin for nodes.

**SwitchData=energy** Specify the data types to be collected by the plugin for switches.

**ColdDoorData=temp** Specify the data types to be collected by the plugin for cold doors.

**MinWatt=<number>** Minimum recorded power consumption, in watts.

**MaxWatt=<number>** Maximum recorded power consumption, in watts.

**MinTemp=<number>** Minimum recorded temperature, in celsius.

**MaxTemp=<number>** Maximum recorded temperature, in celsius.

**EnergyRRA**=<name> Energy RRA name.

**TempRRA=<name>** Temperature RRA name.

**EnergyPathRRD=<path>** Pathname of energy RRD file.

**TempPathRRD=<path>** Pathname of temperature RRD file.

## Example 1 - Node energy monitoring using acct\_gather\_energy/rapl

```
[sulu] (slurm) mnp> scontrol show config
...
AcctGatherEnergyType = acct_gather_energy/rapl
AcctGatherNodeFreq = 30 sec
...

[sulu] (slurm) mnp> scontrol show node n15
NodeName=n15 Arch=x86_64 CoresPerSocket=8
    CPUAlloc=0 CPUErr=0 CPUTot=32 CPULoad=0.00 Features=(null)
    Gres=(null)
    NodeAddr=drak.usrnd.lan NodeHostName=drak.usrnd.lan
    OS=Linux RealMemory=1 AllocMem=0 Sockets=4 Boards=1
    State=IDLE ThreadsPerCore=1 TmpDisk=0 Weight=1
    BootTime=2013-08-28T09:35:47 SlurmdStartTime=2013-09-05T14:31:21
    CurrentWatts=121 LowestJoules=69447 ConsumedJoules=8726863
    ExtSensorsJoules=n/s ExtSensorsWatts=0 ExtSensorsTemp=n/s
```

## Example 2 - Energy accounting using acct\_gather\_energy/rapl

```
[sulu] (slurm) mnp> scontrol show config
JobAcctGatherType
                       = jobacct_gather/linux
JobAcctGatherFrequency = task=10
AcctGatherEnergyType
                       = acct_gather_energy/rapl
                       = accounting_storage/slurmdb
AccountingStorageType
[sulu] (slurm) mnp> srun test/memcputest 100 10000 &
[1] 20712
[sulu] (slurm) mnp> 100 Mb buffer allocated
[sulu] (slurm) mnp> squeue
            JOBID PARTITION
                                NAME
                                        USER ST
                                                       TIME NODES NODELIST(REASON)
              120 drak-only memcpute
                                       slurm R
                                                       0:03
                                                                1 n15
[sulu] (slurm) mnp> sstat -j 120 -o ConsumedEnergy
ConsumedEnergy
-----
         2149
[sulu] (slurm) mnp> sstat -j 120 -o ConsumedEnergy
ConsumedEnergy
-----
         2452
[sulu] (slurm) mnp> sstat -j 120 -o ConsumedEnergy
ConsumedEnergy
-----
         2720
[sulu] (slurm) mnp> Finished: j = 10001, c = 2990739969
[1]+ Done
                            srun test/memcputest 100 10000
[sulu] (slurm) mnp> sacct -j 120 -o ConsumedEnergy
ConsumedEnergy
-----
         3422
```

# **Example 3 - Energy accounting using acct\_gather\_energy/ipmi**

```
[root@cuzco108 bin]# scontrol show config
JobAcctGatherType
                        = jobacct_gather/linux
JobAcctGatherFrequency = task=10
AcctGatherEnergyType
                        = acct_gather_energy/ipmi
AccountingStorageType
                        = accounting_storage/slurmdb
[root@cuzco108 bin]# cat /usr/local/slurm2.6/etc/acct_gather.conf
EnergyIPMIFrequency=10
#EnergyIPMICalcAdjustment=yes
EnergyIPMIPowerSensor=1280
[root@cuzco108 bin]# srun -w cuzco113 memcputest 100 10000 &
[1] 26138
[root@cuzco108 bin]# 100 Mb buffer allocated
[root@cuzco108 bin]# squeue
                                                        TIME NODES NODELIST(REASON)
             JOBID PARTITION
                                 NAME
                                          USER ST
               101 exclusive memcpute
                                                        0:04
                                                                  1 cuzco113
                                          root R
[root@cuzco108 bin]# sstat -j 101 -o ConsumedEnergy
ConsumedEnergy
           570
[root@cuzco108 bin]# sstat -j 101 -o ConsumedEnergy
ConsumedEnergy
         1.74K
```

### **Example 3 - continued**

# **Example 4 - Node energy and temperature monitoring using ext\_sensors/rrd**

```
[root@cuzco0 ~]# scontrol show config
ExtSensorsType
                        = ext_sensors/rrd
ExtSensorsFreq
                       = 10 sec
[root@cuzco108 slurm]# cat /usr/local/slurm2.6/etc/ext_sensors.conf
# External Sensors plugin configuration file
JobData=energy
NodeData=energy, temp
EnergyRRA=1
EnergyPathRRD=/BCM/data/metric/%n/Power_Consumption.rrd
TempRRA=1
TempPathRRD=/BCM/data/metric/%n/Temperature.rrd
MinWatt=4
MaxWatt=200
[root@cuzco0 ~]# scontrol show node cuzco109
NodeName=cuzco109 Arch=x86_64 CoresPerSocket=4
  CPUAlloc=0 CPUErr=0 CPUTot=8 CPULoad=0.00 Features=(null)
   Gres=(null)
   NodeAddr=cuzco109 NodeHostName=cuzco109
   OS=Linux RealMemory=24023 AllocMem=0 Sockets=2 Boards=1
   State=IDLE ThreadsPerCore=1 TmpDisk=0 Weight=1
   BootTime=2013-09-03T17:39:00 SlurmdStartTime=2013-09-10T22:58:10
   CurrentWatts=0 LowestJoules=0 ConsumedJoules=0
   ExtSensorsJoules=4200 ExtSensorsWatts=105 ExtSensorsTemp=66
```

## Example 5 - Energy accounting comparison using ext\_sensors/rrd and acct\_gather\_energy/ipmi

The accuracy/consistency of energy measurements may be inaccurate if the run time of the job is short and allows for only a few samples. This effect should be reduced for longer jobs.

The following example shows that the **ext\_sensors/rrd** and **acct\_gather\_energy/ipmi** plugins produce very similar energy consumption results for a MPI benchmark job using 4 nodes and 32 CPUs, with a run time of ~9 minutes.

#### **Example 5 - continued**

[root@cuzco108 bin]# scontrol show config | grep acct\_gather\_energy

#### acct\_gather\_energy/ipmi

```
AcctGatherEnergyType = acct_gather_energy/ipmi
[root@cuzco108 bin]# srun -n32 --resv-ports ./cg.D.32 &
[root@cuzco108 bin]# squeue
           JOBID PARTITION
                           NAME
                                     USER ST
                                                  TIME NODES NODELIST(REASON)
             122 exclusive cq.D.32
                                      root R
                                                  0:02
                                                           4 cuzco[109,111-113]
[root@cuzco108 bin]# sacct -o "JobID%5, JobName, AllocCPUS, NNodes%3, NodeList%22, State, Start, End, Elapsed, ConsumedEnergy%9"
                                                               Start
        JobName AllocCPUS NNo
                                NodeList State
 127
                      32 4 cuzco[109,111-113] COMPLETED 2013-09-12T23:12:51 2013-09-12T23:22:03 00:09:12 490.60K
       cq.D.32
ext sensors/rrd
[root@cuzco108 bin]# scontrol show config | grep ext_sensors
              = ext_sensors/rrd
ExtSensorsType
[root@cuzco108 bin]# srun -n32 --resv-ports ./cq.D.32 &
[root@cuzco108 bin]# squeue
           JOBID PARTITION
                                      USER ST
                             NAME
                                                  TIME NODES NODELIST(REASON)
             128 exclusive cg.D.32
                                     root R
                                                  0:02
                                                           4 cuzco[109,111-113]
[root@cuzco108 bin]# sacct -o "JobID%5, JobName, AllocCPUS, NNodes%3, NodeList%22, State, Start, End, Elapsed, ConsumedEnergy%9"
JobID
        JobName AllocCPUS NNo
                              NodeList
                                                      State
                                                                        Start
                                                                                                  Elapsed ConsumedE
                      32 4 cuzco[109,111-113] COMPLETED 2013-09-12T23:27:17 2013-09-12T23:36:33 00:09:16 498.67K
 128
       cg.D.32
```

## **Profiling Configuration**

Configuration parameters

The profile plugin is enabled in the **slurm.conf** file, but is internally configured in the **acct\_gather.conf** file.

### slurm.conf parameters

- AcctGatherProfileType=acct\_gather\_profile/hdf5 enables the HDF5 Profile Plugin
- JobAcctGatherFrequency={energy=freq {,lustre=freq {,network=freq, {task=freq}}}} sets default sample frequencies for data types.
- One or more of the following plugins must also be configured.
  - AcctGatherEnergyType=acct\_gather\_energy/ipmi
  - AcctGatherEnergyType=acct\_gather\_energy/rapl
  - AcctGatherFilesystemType=acct\_gather\_filesystem/lustre
  - AcctGatherInfinibandType=acct gather infiniband/ofed
  - JobAcctGatherType=job\_acct\_gather/linux

## Sample conf files

DebugFlags=Profile

#### slurm.conf

```
JobAcctGatherType=jobacct_gather/linux
JobAcctGatherFrequency=energy=5, lustre=60, network=60, task=60
AcctGatherEnergyType=acct_gather_energy/ipmi
AcctGatherFilesystemType=acct_gather_filesystem/lustre
AcctGatherInfinibandType=acct_gather_infiniband/ofed
```

AcctGatherProfileType=acct gather profile/hdf5

### acct\_gather.conf

```
# Parameters for AcctGatherEnergy/ipmi plugin
EnergyIPMIFrequency=10
EnergyIPMICalcAdjustment=yes
#
# Parameters for AcctGatherProfileType/hdf5 plugin
ProfileHDF5Dir=/app/Slurm/profile_data
# Parameters for AcctGatherInfiniband/ofed plugin
InfinibandOFEDFrequency=4
InfinibandOFEDPort=1
```

## **Energy Data**

- AcctGatherEnergyType=acct\_gather\_energy/ipmi is required in slurm.conf to collect energy data.
- JobAcctGatherFrequeny=Energy=<freq> should be set in either slurm.conf or via acctg-freq command line option.

The IPMI energy plugin also needs the EnergyIPMIFrequency value set in the acct\_gather.conf file. This sets the rate at which the plugin samples the external sensors. This value should be the same as the energy=sec in either JobAcctGatherFrequency or --acctg-freq.

Note that the IPMI and profile sampling is not synchronous. The profile sample simply takes the last available IPMI sample value. If the profile energy sample is more frequent than the IPMI sample rate, the IPMI value will be repeated. If the profile energy sample is greater than the IPMI rate, IPMI values will be lost.

Also note that smallest effective IPMI (EnergyIPMIFrequency) sample rate for 2013 era Intel processors is 3 seconds.

Note that Energy data is collected for the entire node so it is only meaningful for exclusive allocations.

Each data sample in the Energy Time Series contains the following data items.

**Date Time** Time of day at which the data sample was taken.

This can be used to correlate activity with other sources such as logs.

**Time** Elapsed time since the beginning of the step.

**Power**Power consumption during the interval.

**CPU Frequency** CPU Frequency at time of sample in kilohertz.

### **Lustre Data**

- AcctGatherFilesystemType=acct\_gather\_filesystem/lustre is required in Slurm.conf to collect lustre data.
- JobAcctGatherFrequeny=Lustre=<freq> should be set in either Slurm.conf or via -acctg-freq command line option.
- Each data sample in the Lustre Time Series contains the following data items.

**Date Time** Time of day at which the data sample was taken.

This can be used to correlate activity with other sources such as logs.

**Time** Elapsed time since the beginning of the step.

**Reads** Number of read operations.

MegabytesRead Number of megabytes read.

**Writes** Number of write operations.

**MegabytesWrite** Number of megabytes written.

### **Network (Infiniband) Data**

- AcctGathertInfinibandType=acct\_gather\_infiniband/ofed is required in Slurm.conf to collect Network data.
- JobAcctGatherFrequeny=Network=<freq> should be set in either Slurm.conf or via -acctg-freq command line option.
- Each data sample in the Network Time Series contains the following data items.

**Date Time** Time of day at which the data sample was taken.

This can be used to correlate activity with other sources such as logs.

**Time** Elapsed time since the beginning of the step.

**PacketsIn** Number of packets coming in.

MegabytesIn Number of megabytes coming in through the interface.

**PacketsOut** Number of packets going out.

**MegabytesOut** Number of megabytes going out through the interface.

### **Task Data**

- JobAcctGatherType=jobacct\_gather/linux is required in Slurm.conf to collect task data
- JobAcctGatherFrequeny=Task=<freq> should be set in either Slurm.conf or via -acctg-freq command line option.

The frequency should be set to at least 30 seconds for CPU utilization to be meaningful (since the resolution of cpu time in linux is 1 second)

• Each data sample in the Task Time Series contains the following data items.

**Date Time** Time of day at which the data sample was taken.

This can be used to correlate activity with other sources such as logs.

**Time** Elapsed time since the beginning of the step.

**CPUFrequency** CPU Frequency at time of sample.

**CPUTime**Seconds of CPU time used during the sample.

**CPUUtilization** CPU Utilization during the interval.

**RSS** Value of RSS at time of sample.

**VMSize** Value of VM Size at time of sample.

**Pages** Pages used in sample.

**ReadMegabytes** Number of megabytes read from local disk.

WriteMegabytes Number of megabytes written to local disk.

### **Emulation and Performance Evaluation**

### Activating emulation technique within SLURM

- •Multiple slurmd technique can be used to experiment with larger scales:
- the idea is that multiple slurmd deamons use the same IP address but different ports
  - all controller side plugins and mechanisms will function
  - ideal for scheduling, internal communications and scalability experiments

- 1. You need to run ./configure with -enable-multiple-slurmd parameter (make, make install, etc)
- 2. Perform the necessary changes in the slurm.conf file similarly the following example:

### Activating emulation technique within SLURM

```
SlurmdPidFile=/usr/local/slurm-test/var/run/slurmd-%n.pid
SlurmdSpoolDir=/tmp/slurm-%n
SlurmdLogFile=/tmp/slurmd-%n.log
FastSchedule=2
PartitionName=exclusive Nodes=virtual[0-40] Default=YES MaxTime=INFINITE State=UP Priority=10
Shared=EXCLUSIVE
NodeName=DEFAULT Sockets=2 CoresPerSocket=8 ThreadsPerCore=1 RealMemory=21384 State=IDLE
NodeName=virtual0 NodeHostName=nazgul NodeAddr=127.0.0.1 Port=17000.
NodeName=virtual1 NodeHostName=nazgul NodeAddr=127.0.0.1 Port=17001
NodeName=virtual2 NodeHostName=nazgul NodeAddr=127.0.0.1 Port=17002
......
```

#### 3. You can start the slurmd deamons with:

Either through a script such as:

```
for i in {0..40}; do slurmd -N virtual$i; done
```

Or by exporting: MULTIPLE\_SLURMD="\$(grep NodeHostName=\$(hostname) / etc/slurm.conf | cut -d ' ' -f 1 | cut -d'=' -f 2)"

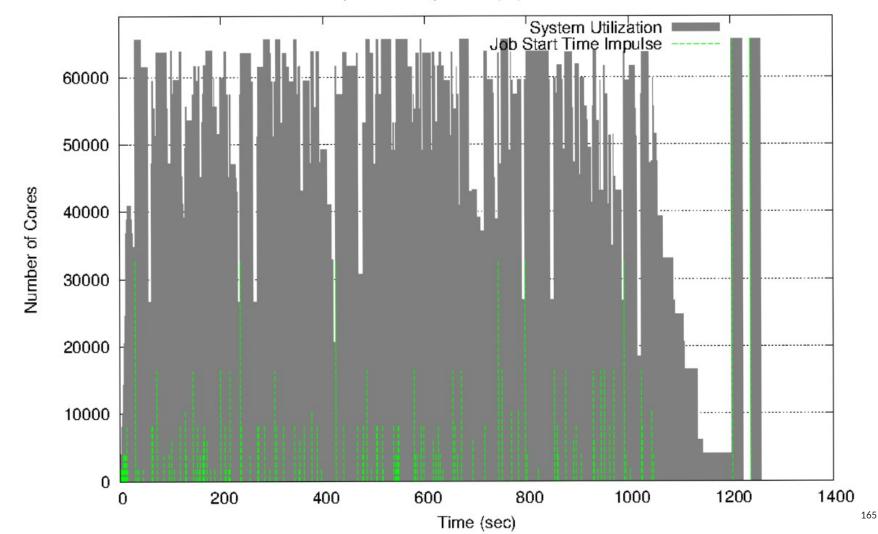
on /etc/sysconfig/slurm and starting with /etc/init.d/slurm

### Examples of performance evaluation with emulation

### 4096 emulated nodes upon 400 physical nodes

© Bull,

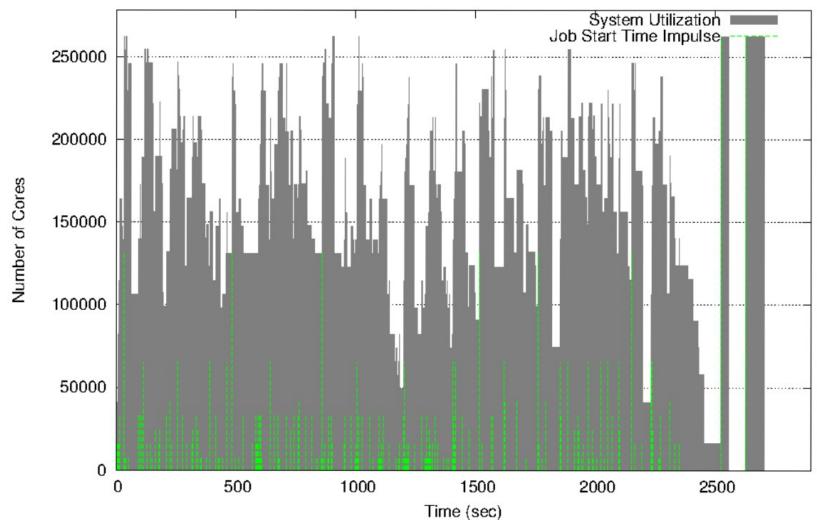
System utilization for Light ESP synthetic workload of 230jobs and SLURM upon 4096 nodes (16cpu/node) cluster (emulation upon 400 physical nodes)



### Examples of performance evaluation with emulation

### 16384 emulated nodes upon 400 physical nodes

System utilization for Light ESP synthetic workload of 230jobs and SLURM upon 16384 nodes cluster (emulation upon 400 physical nodes)



- Introduction
- SLURM scalable and flexible RJMS
- Part 1: Basics
  - Overview, Architecture, Configuration files, Partitions, Plugins, Reservations, Reconfiguration
- Part 2: Advanced Configuration
  - Accounting, Scheduling, Allocation, Network Topology Placement, Generic Resources, Licenses Management, Energy Reduction Techniques
- Part 3: Experts Configuration
  - CPU Management, Isolation with cgroups, Power Management, Simulation and evaluation
- Upcoming Features



### New Slurm features under development

- Heterogeneous Environment
  - Asymmetric Resources and MPMD model
  - GPU Affinity
- Scalability
  - Support of PMI-x project
  - Messages Aggregation
  - HDF5 Profiling Framework
- Power Management and Energy Efficiency
  - Extension of Energy Accounting and Power Profiling Framework
  - Power-Capping logic in Job Scheduling
  - Energetic Fairsharing

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