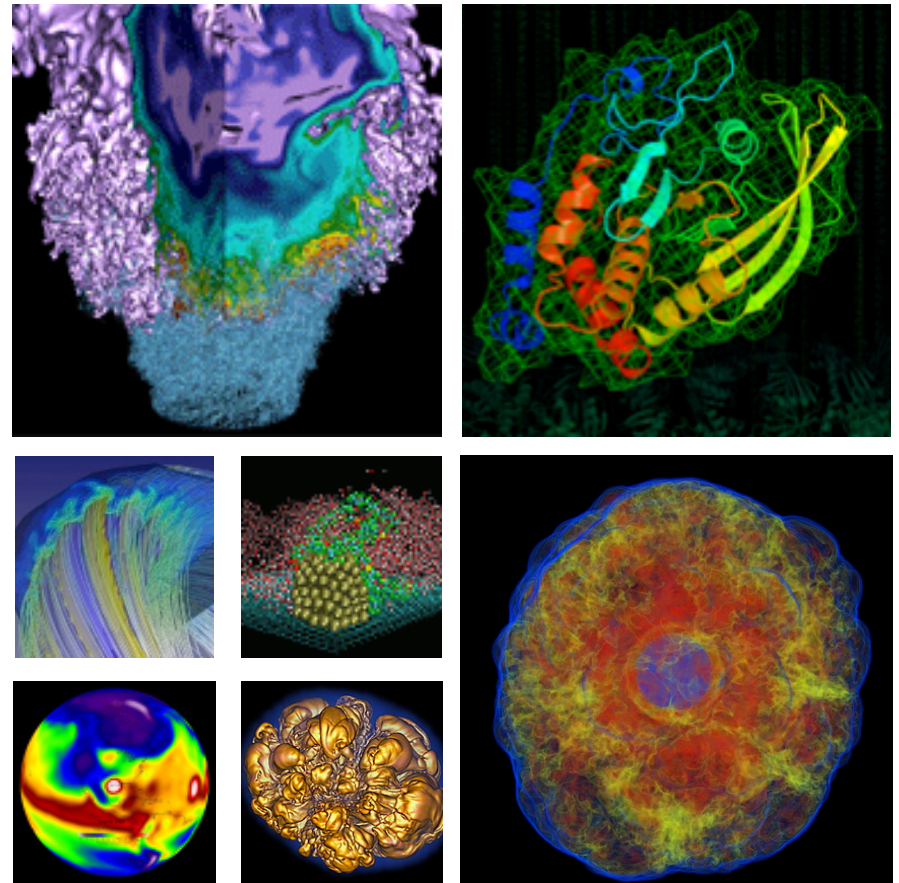


Getting Started at NERSC: Running Jobs



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Jobs at NERSC



- **Most jobs are parallel, using 10s to 100,000+ cores**
- **Production runs execute in batch mode**
- **Interactive and debug jobs are supported for up to 30 minutes**
- **Typically run times are a few to 10s of hours.**
 - Each machine has different limits.
 - Limits are necessary because of MTBF and the need to accommodate 5,500 users' jobs
- **Also a number of 'serial' jobs**
 - Typically some kind of pleasantly parallel simulation

What system is best for your job?

Large-Scale Computing Systems

Edison (NERSC-7): Cray XC30

- 5,192 compute nodes, 124,608 cores
- 236 Tflop/s on applications, 2.39 Pflop/s peak

Hopper (NERSC-6): Cray XE6

- 6,384 compute nodes, 153,216 cores
- 144 Tflop/s on applications; 1.3 Pflop/s peak



Midrange

Thousands of cores

Carver

- IBM iDataplex cluster, 1,202 nodes
- 9984 cores; 106 TF



PDSF (HEP/NP)

- ~2K core cluster

GenePool (JGI)

- ~5K core cluster
- 2.1 PB Isilon File System

NERSC Global Filesystem (NGF)

Uses IBM's GPFS

- 12.7 PB capacity
- 150 GB/s of bandwidth



HPSS Archival Storage

- 240 PB capacity
- 5 Tape libraries
- 200 TB disk cache



Analytics & Testbeds



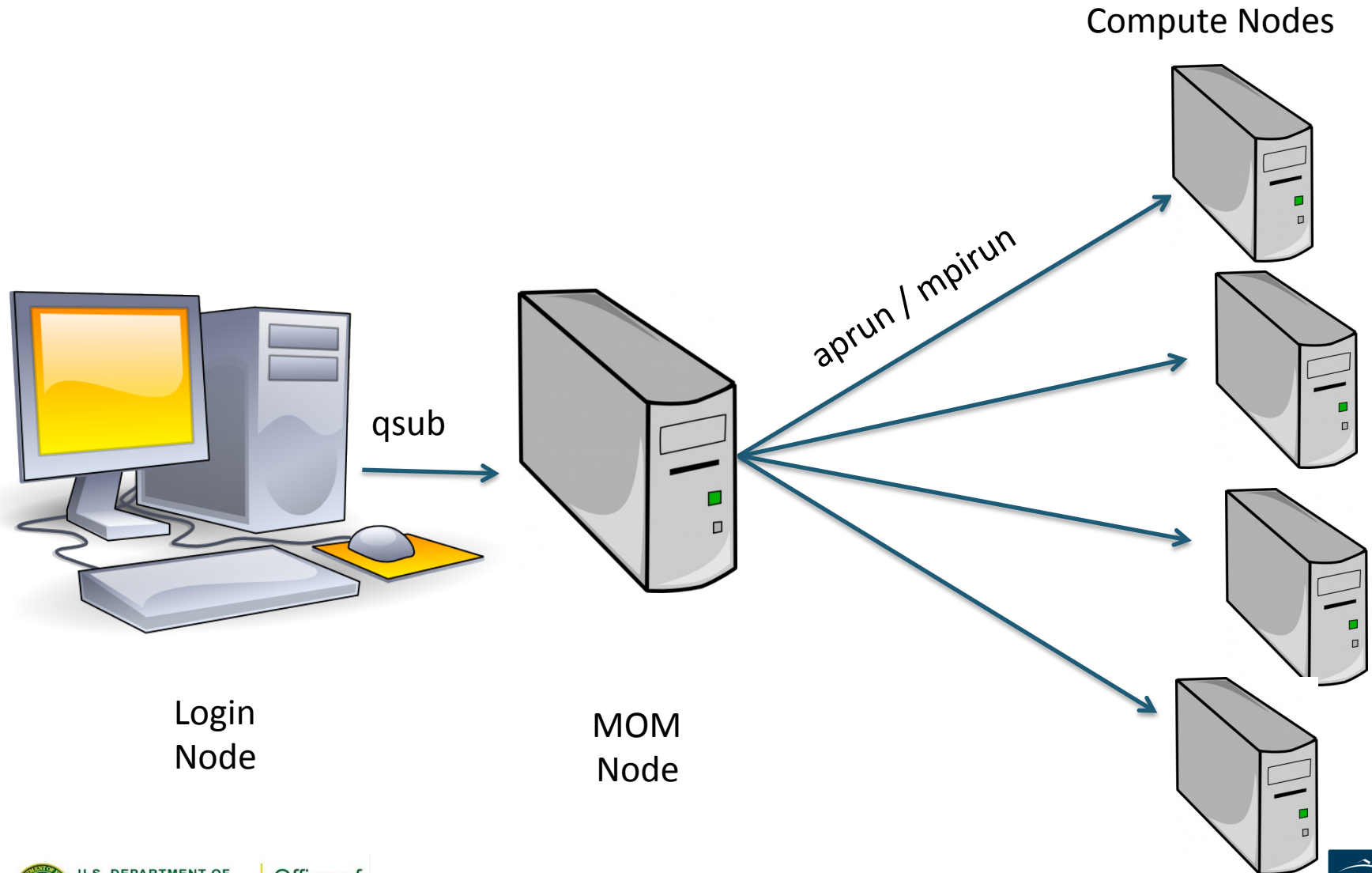
Dirac 48 Fermi GPU nodes

Login Nodes and Compute Nodes



- Each supercomputer has 3 types of nodes that you will use
 - Login nodes
 - Compute nodes
 - Job launcher or “MOM” nodes
- Login nodes
 - Edit files, compile codes, run UNIX commands
 - Submit batch jobs
 - Run short, small utilities and applications
- Compute nodes
 - Execute your application; dedicated to your job
 - No direct login access
- Job launcher or “MOM” nodes
 - Execute your batch script commands
 - Carver: “head” compute node; Edison / Hopper: shared “service” node; not a compute node

Launching Parallel Jobs



Launching Parallel Jobs

- A “job launcher” executes your code
 - Distributes your executables to all your nodes
 - Starts concurrent execution of N instances of your program
 - Manages execution of your application
 - On Edison / Hopper: the job launcher is called “aprun”
 - On Carver: “mpirun”
- Only the job launcher can start your job on compute nodes
- You can’t run the job launcher from login nodes

Submitting Batch Jobs

- To run a job on the compute nodes you must write a “batch script” that contains
 - Batch directives to allow the system to schedule your job
 - An `aprun` or `mpirun` command that launches your parallel executable
- Submit the job to the queuing system with the `qsub` command
 - `%qsub my_batch_script`

Edison - Cray XC30



- 124,608 cores, 5,192 nodes
- “Aries” interconnect
- 2 12-core Intel ‘Ivy Bridge’ 2.4 GHz processors per node
- 24 processor cores per node, 48 with hyperthreading
- 64 GB of memory per node
- 332 TB of aggregate memory
- 2.7 GB memory / core for applications
- /scratch disk quota of 10 TB
- 7.6 PB of /scratch disk
- Choice of full Linux operating system or optimized Linux OS (Cray Linux)
- Intel, Cray, and GNU compilers

- **Login Nodes**

- Edison's login nodes run a full Linux operating system and provide support services for the system. When you connect to Edison with SSH, you land on the login nodes. These nodes are shared by many users; please do not run applications on the login nodes.

- **Compute Nodes**

- The 5,192 compute nodes are dedicated to running scientific applications. A job is given exclusive access to each node it requests for the entirety of the job's run time. Since Edison has 24 cores on each node, the minimum number of cores per job is 24.

- **Job Host (MOM) Nodes**

- MOM nodes are servers that execute batch job commands. These nodes are shared by many users and thus are not intended for compute- or memory-intensive applications.

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 ./my_executable
```

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 ./my_executable
```

Job directives: instructions for the batch system

- Submission queue
- How many nodes to reserve for your job
- How long to reserve those nodes
- Optional: what to name STDOUT files, what account to charge, whether to notify you by email when your job finishes, etc.

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 ./my_executable
```

Change from home directory to job submission directory

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 ./my_executable
```

Launches parallel executable on the compute nodes.
Carries over (partial) login environment and controls how your executable accesses the memory on each processor.

Sample Hopper Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 ./my_executable
```

mppwidth is number of compute cores requested (/24 for number of nodes).
Must match with number of tasks requested (-n)

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=192
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 -N 12 ./my_executable
```

-N = number of tasks per node

Might do this to get more memory / task

Hybrid OpenMP/MPI

```
#PBS -q regular
#PBS -l mppwidth=96
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
export OMP_NUM_THREADS=6
aprun -n 16 -d 6 -N 4 -S 1 ./hybrid.x
```

16 tasks (-n), 4 on each node (-N), 6 OpenMP threads / task (-d), separate tasks onto individual NUMA nodes (-S)

Many more examples on www.nersc.gov

Sample Edison Batch Script - MPI

```
#PBS -q debug
#PBS -l mppwidth=48
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
aprun -n 96 -j 2 ./my_executable
```

Turn on hyperthreading (-j)

Carver - IBM iDataPlex

- 9,984 compute cores
- 1,202 compute nodes
- 2 quad-core Intel Nehalem 2.67 GHz processors per node
- 8 processor cores per node
- 24 GB of memory per node (48 GB on 80 "fat" nodes)
- 2.5 GB / core for applications (5.5 GB / core on "fat" nodes)
- InfiniBand 4X QDR



- NERSC global /scratch directory quota of 20 TB
- Full Linux operating system
- PGI, GNU, Intel compilers

Sample Carver Batch Script - MPI

```
#PBS -q debug
#PBS -l nodes=16:ppn=8
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
mpirun -n 128 ./myexecutable
```

↑
MPI tasks

Interactive Parallel Jobs

- You can run small parallel jobs interactively for up to 30 minutes (ex. is for Hopper)

```
% qsub -I -l mppwidth=48
```

```
[wait for job to start]
```

```
% cd $PBS_O_WORKDIR
```

```
% aprun -n 48 ./mycode.x
```

- Carver has a special queue for running serial jobs
 - A single process running on a single node
 - Each serial node can run up to 12 jobs from different users

```
#PBS -q serial
#PBS -l walltime=00:10:00
#PBS -N my_job

cd $PBS_O_WORKDIR
./myexecutable
```

Monitoring Your Job



- Once your job is submitted, it will start when resources are available

- Monitor it with

- `qstat -a`
- `qstat -u username`
- `showq`
- `qs`
- NERSC web site “Queue Look”

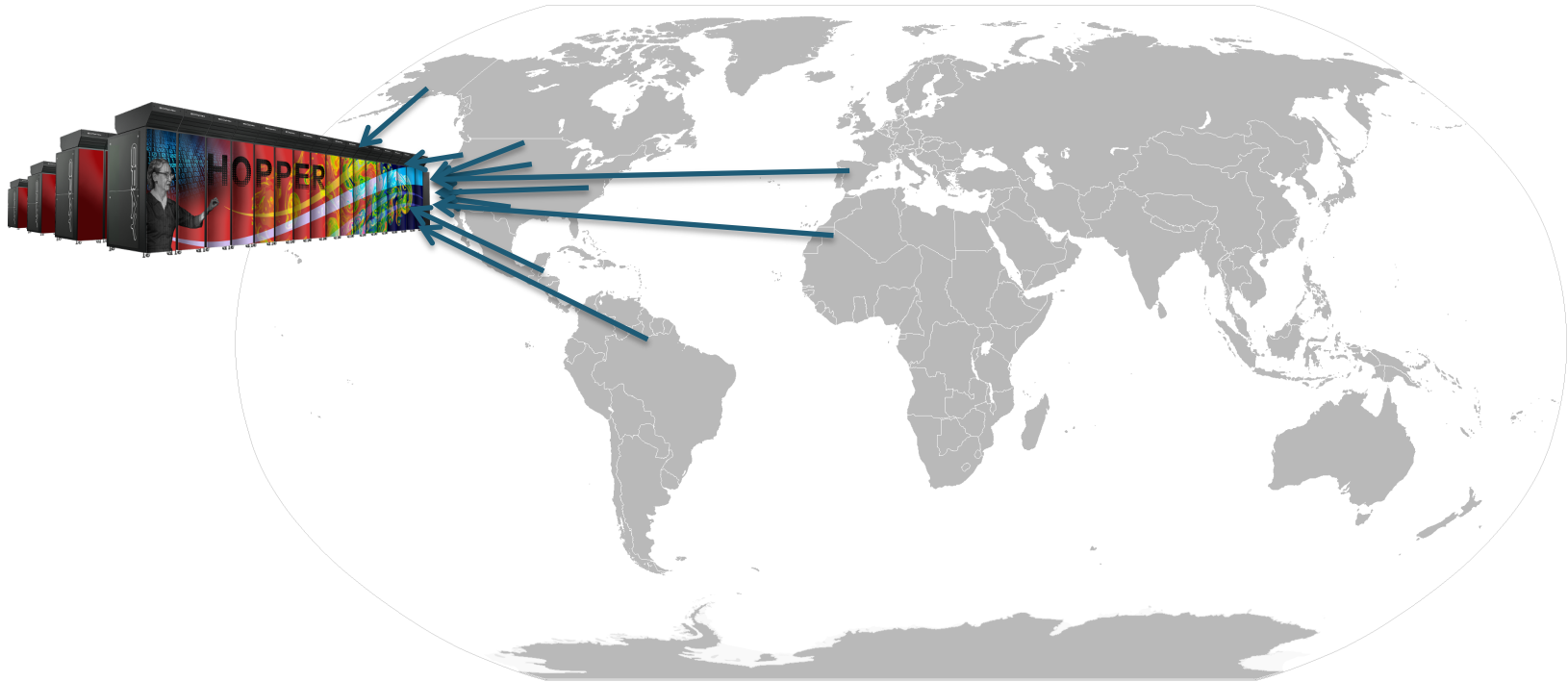
<https://www.nersc.gov/users/live-status/global-queue-look/>

- NERSC web site “Completed Jobs”

<https://www.nersc.gov/users/job-logs-and-analytics/completed-jobs/>

I qsub'd my job, but it's not running!

- You are not alone!



How do we handle the demand?

That's what the batch queue is for!



source: itu.dk

Your jobs will wait until the resources are available for them to run

Your job's place in the queue is a mix of time and priority, so line jumping is allowed, but it may cost more

Job Limits

There are per user, per machine job limits. See the NERSC web site for details. Here are the limits on Edison as of Jan. 28, 2014.

Specify these queues
(with #PBS -q queue_name)

NEVER these!

Submit Queue	Execution Queue	Nodes	Physical Cores	Max Wallclock (hours)	Relative Priority	Run Limit	Queued Limit	Queue Charge Factor
debug	debug	1-512	1-12,288	30 mins	2	2	2	1
ccm_int ¹	ccm_int	1-512	1-12,288	30 mins	2	2	2	1
regular	reg_1hour	1-256	1-6144	1 hr	3	8	8	1
	reg_small	1-682	1-16,368	48 hrs	3	16	16	1
	reg_med	683-2048	16,369-49,152	36 hrs	3	4	4	0.75
	reg_big	2049-4096	49,153-98,304	36 hrs	2	2	2	0.75
	reg_xbig	4097-5000	98,305-120,000	12 hrs	2	2	2	0.75
ccm_queue	ccm_queue	1-682	1-16,368	48 hrs	3	16	16	1
premium	premium	1-2048	1-49,152	12	1	1	1	2
low	low	1-682	1-16,368	24	4	8	8	0.5
killable ²	killable	1-682	1-16,368	48 hrs	3	8	8	1

- Submit shorter jobs.
 - Checkpoint if possible
 - Take advantage of ‘backfill’ opportunities
 - Run jobs just before maintenance
- Make sure the wall clock time you request is accurate. As noted above, shorter jobs are easier to schedule. Many users unnecessarily enter the largest wall clock time possible as a default.

NERSC 40 YEARS
at the
FOREFRONT

	Hours Requested																																																		
Nodes	<1	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	25	26	27	28	29	30	31	32	33	34	35	36	37	38	39	40	41	42	43	44	45	46	47	48	48+	
1	1.6	5.3	4.1	5.6	5.8	13	38	58		24	9.8	17	18	110	67	15	31	16	42	6.2	17	10	26	26	29	11	19	20	57	26	42	13	12	27	34	21	46		18	26	26	17	20		28	20	21	24	43	29	
2	1.7	7.5	8.0	9.6	22	15	25	8.6	21		14	15	13	23	18	22	23	15	14	30	34	34	31	25	30	41	21	35	33	20	19	25		41		36	29	47	40	0.1	33	45		35	17	58	31	33	33	58	76
3	1.4		17.4	8.5	1.4	4.8	40	48	22	11	-2	590.3	155	0	19	28	29	19	26	14	23	23	51		35	37	45	66	89		32	40	51		42		50	112	44	31		35			8.3	22	39	37	93	24	
4	3.6		11.6	3.1	14	10	0	12	14	99	16		27	27	29	38	16	24	26	39	20	71	28	48		14	29	52	60	44		35		112		54		69	54	25	65	148					93	65	44		
5	2.2		1.1	7.2	6.0	5.9	9.7	20	13	16		6.5	49	8.6	19	13	38	25	12	28	35	33	58	43	25	50	76	61			23	43		47			43	49		44		44		45				25	74	32	
6	1.1		6.0	3.5	10	6.4	13	11	13	41		11	40	21	19	24	21	23	34	44	5	32		43		35	84	80		27		78		58	34	56	49	28	53	59		42		53			56	93	102	34	
7	0.4		5.4	2.5	7.5	9.6	6.5	14	3.2	15		7.8	38	31	24							47			83	16	57					37		47			55	19						104			64	35			
8	1.8		16	8.4	36	9.8	16	19	17	16		16	59	41	22	4.7	27	36	37			33	57	64			73	68	18	18		56		25		75	44	28						86	17		77	67	35		
9	0.4		2.1	2.5	5.8	8.7	15	11	12	14			32	18	18			36	23								45					236							63			51		90	49	72	59	104	16		
10	1.8		7.2	8.6	4.8	13	11	12	17	21		18	28	26	23	50	81	72	40			13	25	80		72	50	110	60			62										55		113	91		57	87	76		
11	10		26	7.0	6.1	16	13	30	15	15		11	22	27	31	38	27	31	25	39	45	61	75			125	56		41				59										52		48	58	81	102	45		
12	2.4		9.6	10.5	3	23	47	96	16	36		31	254	18	32	9.5		45	40		22	53	60			113	57	61		34		36					138			38	63		68	37	63	47	59	42			
13	0.4		1.4	7.5	10	19	136	4	8		10	80		18								48						57				50					15					64		70		66					
14	1.6		1.1	18	12	8.7	22	17	14	27		27	27		29					105		97				37	54				38											53		151			109	120			
15	12		6.7	15	16	7.7	9.9	62	81	14		47	53	37	25	47	86	38	17	38		72			65	204	54		48			44				1	93						51			77	87				
16	1.6		1.9	10	27	11	22	39	18	18		30	27	28	22	124	42	40	43	29	66	58	77		65	65	79	33	45		96	68	140		53						105				80	55	32				
17-19	1.3		26	19	5.3	10	8	0	13	46		21	30	11	43	104	27	30		75		76			144	112	132			152	63	36		116			64	98		49		69		67			77	53			
20-23	8.9		8.3	9.3	19	20	19	23	24	29		46	32	25	29	68	53	57	154			67	57	76	41	61	66	59			39	75		202			59	61			65		27		44	41	75	36			
24-31	2.6		2.6	4.8	11	15	24	48	15	46		30	19	11	25	166	47	43	47	69	57	67	53		28	66	125	78			86	39	66									94			100	40	40	87	104		
32-47	1.5		9.2	9.5	26	18	27	58	18	36		51	45	28	37	82	51	54	49	65	86	44	60	128	47	66	101	46				21	89		134	43		67	67			84	49	88	47		62	65	79	88	55
48-63	6.0		12	9.3	15	18	18	32	54	83		43	31	65	40		47	64	32		74	74	52			71	114	62			86		64		69		108	72		34		58		106	95	77	80	98			
64-127	1.6		23	20	21	31	29	36	45	53		33	39	43	52	77	67	59	73	60	86	61	72	45	86	94	78	63	61		105	119	59		64		127	64			42	65			59	58		70	101		
128-255	4.8	-178.1	21	38	40	50	54	45	77		51	33	149	48	66	130	58	73	73	78		63		55	79	101					41	73		34			175	64				45	50		98		109	102	91	63	
256-511	7.8		41	47	38	83	47	131	30	67		51	56	32	44	43	71	51	79	2.5	105	58	95			98	86	349															68				112	105			
512-1023	12		53	100	91	48	48	65	30	50		94	44	60	56	62	28	66	54	44	36	41	53		82	32	42	16		32	42		87		144		36	82	76							128					
1024-1535	21		31	32	26	301	29	77	35	56		55	90	39	89	39	41		33	65	93		45									72															139	208			
1536-2047	24		43	21	11	35	32	66		48		47		35			39		31		142		19		64																					110					
2048-3071	30		32	59	41	32	45	50	49	45		106		21	46		70	203	78	144	101	33			67	53	31		56																						
3072-4095	25		57	102							130			193								72																													
4096-6143	133		75	75	50	50	0.0	129	60			14																																							
6144-9531	0.0		55		29																																														

How Your Jobs Are Charged

- Your repository account is charged for **each node** your job was **allocated** for the **entire duration** of your job.
 - The minimum allocatable unit is a **node**. Hopper and Edison have 24 cores/node, so your minimum charge is $24 * \text{walltime}$.
- $$\text{MPP hours} = (\# \text{ nodes}) * (\# \text{ cores / node}) * (\text{walltime}) * (\text{QCF}) * (\text{MCF})$$
- mppwidth = 96 for 1 hour. $\text{MPP hours} = (4) * (24) * (1 \text{ hour}) * (1) * (1) * \text{MPP hours charged} = 96 \text{ hours}$
 - Serial jobs on Carver are charged with $\text{walltime} * \text{MCF}$
- If you have access to multiple repos, pick which one to charge in your batch script
 - `#PBS -A repo_name`

Charge Factors & Discounts

- Each machine has a “machine charge factor” (mcf) that multiplies the “raw hours” used
 - Edison has MCF = 2
 - Hopper has MCF = 1
 - Carver has MCF = 1.5
- Queues have “priority charge factors” (pcf) and corresponding relative scheduling priorities
 - Premium pcf = 2.0
 - Low pcf = 0.5
 - Everything else pcf = 1.0
- On Edison:
 - reg_med, reg_big, reg_xbig jobs get a 25% discount

More Information



NERSC Web pages

Hopper:

<http://www.nersc.gov/users/computational-systems/hopper/running-jobs/>

Edison:

<http://www.nersc.gov/users/computational-systems/edison/running-jobs/>

Carver:

<http://www.nersc.gov/users/computational-systems/carver/running-jobs/>

Contact NERSC Consulting:

- Toll-free 800-666-3772
- 510-486-8611, #3
- Email consult@nersc.gov.



Thank You

Sample Carver Batch Script - OpenMP

```
#PBS -q debug
#PBS -l nodes=1:ppn=8
#PBS -l walltime=00:10:00
#PBS -N my_job
```

```
export OMP_NUM_THREADS=8
cd $PBS_O_WORKDIR
./myexecutable
```

Number of threads