



## Profiling Applications to Choose the Right Computing Infrastructure plus Batch Management with HTCondor

#### DOSAR

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### https://osg-htc.org/dosar/ASP2024/ASP2024\_Materials/







- It's okay to move ahead on exercises if you have time
- It's okay to take longer on them if you need to
- If you move along quickly, try the "On Your Own" sections and "Challenges"









 Please ask us questions! ...during the lectures ...during the exercises ...during the breaks ...during the meals ...over dinner ...via email after we depart (see below)

• If I don't know, I'll find the right person to answer your question.







- Profiling your application
- Picking the appropriate resources
- Understand the basics of HTCondor

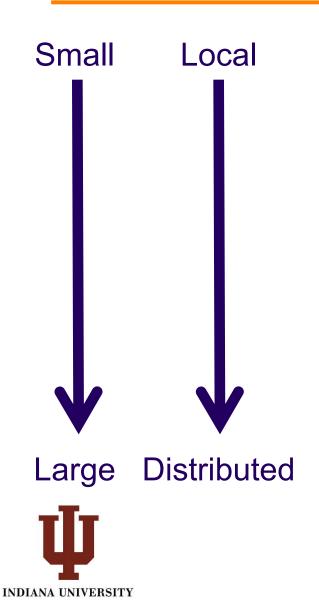






## Let's take one step at a time





- Can you run one job on one computer?
- Can you run one job on another computer?
- Can you run 10 jobs on a set of computers?
- Can you run a multiple job workflow?
- How do we put this all together?

This is the path we'll take in the school



INDIANA UNIVERSITY





- A "headless job"
  - Not interactive/no GUI: how could you interact with 1000 simultaneous jobs?
- A set of input files
- A set of output files
- A set of parameters (command-line arguments)
- Requirements:
  - Ex: My job requires at least 2GB of RAM
  - Ex: My job requires Linux
- Control/Policy:
  - Ex: Send me email when the job is done
  - Ex: Job 2 is more important than Job 1
  - Ex: Kill my job if it runs for more than 6 hours





- Methods to:
  - Submit/Cancel job
  - Check on state of job
  - Check on state of available computers
- Processes to:
  - Reliably track set of submitted jobs
  - Reliably track set of available computers
  - Decide which job runs on which computer
  - Manage a single computer
  - Start up a single job





- Let's assume you have a 'large job' – What factors could make it large?
- Large Data Input or Output or both
- Needs to do heavy calculation
- Needs a lot of memory
- Needs to communicate with other jobs (whether required or not)
- Reads and writes a lot of data/files
- Heavy graphics processing
- Any combination of any of the above







- But some solutions are more "Open" than others.
  - Local Laptop/Desktop
  - Local Cluster
  - HPC System
  - Shared HTC Resources
  - Clouds





# Why is HTC hard?



- The HTC system has to keep track of:
  - Individual tasks (a.k.a. jobs) & their inputs
  - Computers that are available
- The system has to recover from failures
  - There will be failures! Distributed computers means more chances for failures.
- You have to share computers
  - Sharing can be within an organization, or between orgs
  - So you have to worry about security
  - And you have to worry about policies on how you share
- If you use a lot of computers, you have to handle variety:
  - Different kinds of computers (arch, OS, speed, etc..)
  - Different kinds of storage (access methodology, size, speed, etc...)
  - Different networks interacting (network problems are hard to debug!)







- Surprise! HTCondor does this (and more)
- Methods to:
  - Submit/Cancel job. condor\_submit/condor\_rm
  - Check on state of job. <a href="mailto:condor\_q">condor\_q</a>
  - Check on state of avail. computers. condor\_status
- Processes to:
  - Reliably track set of submitted jobs. schedd
  - Reliably track set of avail. computers. collector
  - Decide which job runs on where. negotiator
  - Manage a single computer startd
  - Start up a single job starter







- You can use other systems:
  - PBS/Torque
  - Oracle Grid Engine (né Sun Grid Engine)
  - LSF
  - SLURM
  - ..
- But I won't cover them.
  - My experience is with Condor
  - My bias is with Condor
  - Overlays exist
- What should you learn at the school?
  - How do you think about Computing Resources?
  - How can you do your science with HTC?
  - For now, learn it with Condor, but you can apply it to other systems.





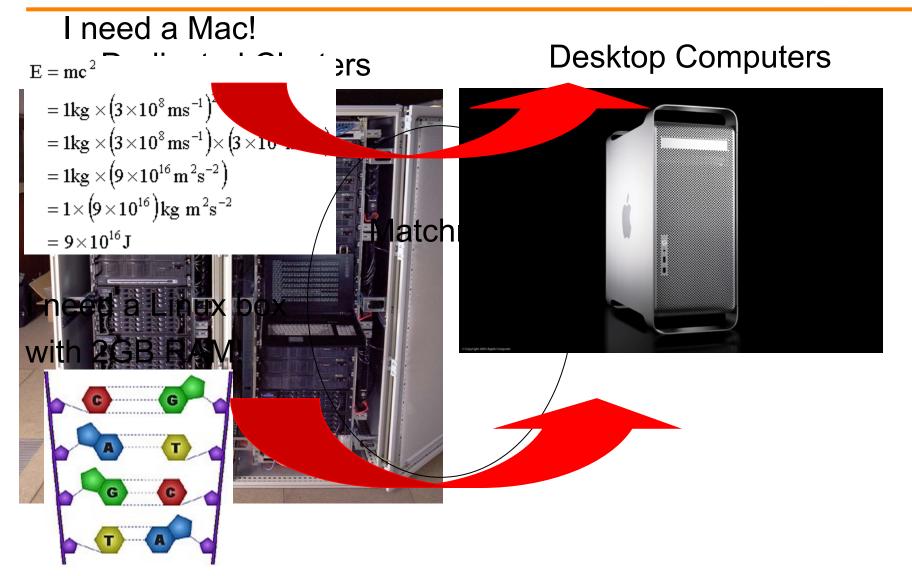




- Please note, we will only scratch the surface of Condor:
  - We won't cover MPI, Master-Worker, advanced policies, site administration, security mechanisms, submission to other batch systems, virtual machines, cron, high-availability, computing on demand, containers.



# Open Science And Matches Computers...



U.S.





- Cluster: A dedicated set of computers not for interactive use
- Pool: A collection of computers used by Condor
  - May be dedicated
  - May be interactive
- Remember:
  - Condor can manage a cluster in a machine room
  - Condor can use desktop computers
  - Condor can access remote computers
  - HTC uses all available resources







- Matchmaking is fundamental to Condor
- Matchmaking is two-way
  - Job describes what it requires:

I need Linux && 8 GB of RAM

– Machine describes what it requires:

I will only run jobs from the Physics department

- Matchmaking allows preferences
  - I need Linux, and I prefer machines with more memory but will run on any machine you provide me

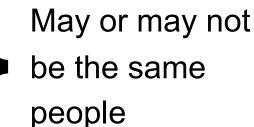








- Condor conceptually divides people into three groups:
  - Job submitters
  - Computer owners
  - Pool (cluster) administrator



All three of these groups have preferences









- ClassAds state facts
  - My job's executable is analysis.exe
  - My machine's load average is 5.6
- ClassAds state preferences

   I require a computer with Linux
- ClassAds are extensible

   They say whatever you
   want them to say











МуТуре	= "Job" ← String
TargetType	= "Machine"
ClusterId	= 1377 <b>← Number</b>
Owner	= "roy"
Cmd	= "analysis.exe"
Requirements = Expression	
(Arch ==	"INTEL")
&& (OpSys ==	"LINUX")
&& (Disk >=	DiskUsage)
&& ((Memory	* 1024)>=ImageSize)









- Condor imposes some schema

   Owner is a string, ClusterID is a number...
- But users can extend it however they like, for jobs or machines
  - AnalysisJobType = "simulation"
  - HasJava\_1\_6 = TRUE
  - ShoeLength = 10
- Matchmaking can use these attributes



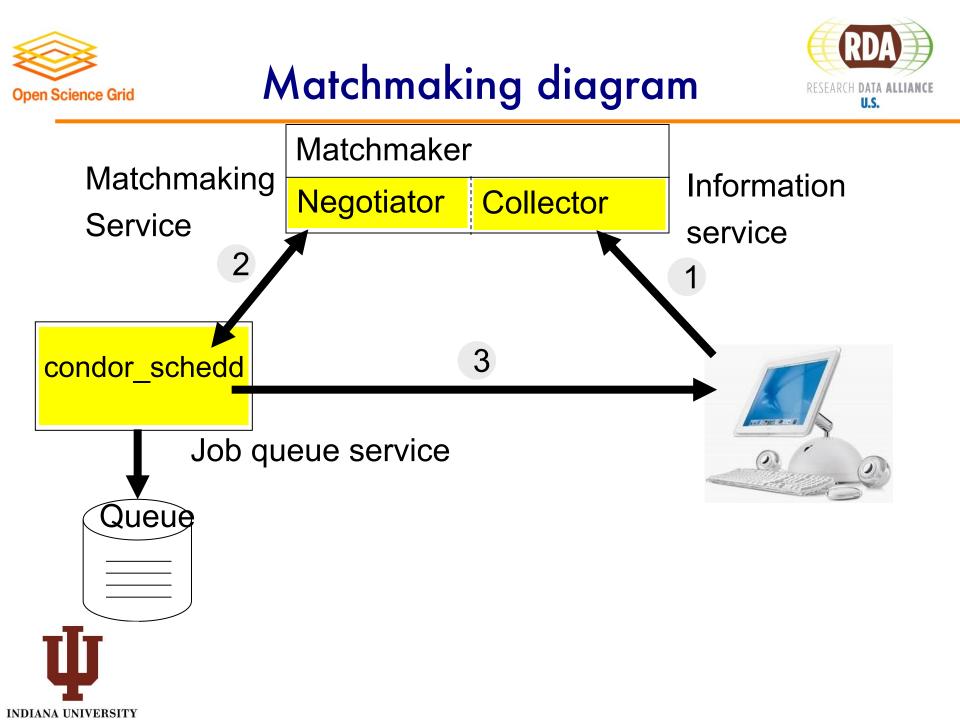






- You won't write ClassAds (usually)
  - You' Il create a simple submit file
  - Condor will write the ClassAd
  - You can extend the ClassAd if you want to
- You won't write requirements (usually)
  - Condor writes them for you
  - You can extend them
  - In some environments you provide attributes instead of requirements expressions











- The computer running the job fails

   Or the network, or the disk, or the OS, or...
- Your job might be *preempted*:
  - Condor decides your job is less important than another, so your job is stopped and another started.
  - Not a "failure" per se, but it may feel like it to you.









- When a job fails or is preempted:
  - It stays in the queue (on the schedd)
  - A note is written to the job log file
  - It reverts to "idle" state
  - It is eligible to be matched again
- Relax! Condor will run your job again





Access to data in Condor



- Option #1: Shared filesystem
  - Simple to use, but make sure your filesystem can handle the load
- Option #2: Condor's file transfer
  - Can automatically send back changed files
  - Atomic transfer of multiple files
  - Can be encrypted over the wire
  - Most common for small applications and data
- Option #3: Remote I/O



# **Condor File Transfer**



- ShouldTransferFiles = YES
  - Always transfer files to execution site
- ShouldTransferFiles = NO
  - Rely on a shared filesystem
- ShouldTransferFiles = IF\_NEEDED
  - Will automatically transfer the files if needed

```
Universe = vanilla
Executable = my_job
Log = my_job.log
ShouldTransferFiles = YES
Transfer_input_files = dataset$(Process), common.data
Queue 600
```



## Transfer\_input\_files can be a URL For example:

transfer\_input\_files = http://www.example.com/input.data







- One submit file can describe lots of jobs
  - All the jobs in a submit file are a *cluster* of jobs
  - Yeah, same term as a cluster of computers
- Each cluster has a unique "cluster number"
- Each job in a cluster is called a "process"
- A Condor "job ID" is the cluster number, a period, and the process number ("20.1")
- A cluster is allowed to have one or more processes.
  - There is always a cluster for every job





The \$(Process) macro



- The initial directory for each job can be specified as run\_\$(Process), and instead of submitting a single job, we use "Queue 600" to submit 600 jobs at once
- The \$(Process) macro will be expanded to the process number for each job in the cluster (0 599), so we'll have "run\_0", "run\_1", ... "run\_599" directories
- All the input/output files will be in different directories!









# Example condor\_submit input file that defines # a cluster of 600 jobs with different directories Universe = vanilla Executable = my\_job Log = my\_job.log Arguments = -arg1 -arg2 Input = my\_job.stdin Output = my\_job.stdout Error = my\_job.stderr InitialDir = run\_\$(Process) Queue 600









### • You can use \$(Process) anywhere:

Universe = vanilla Executable = my\_job Log = my\_job.\$(Process).log Arguments = -randomseed \$(Process) Input = my\_job.stdin Output = my\_job.stdout Error = my\_job.stderr InitialDir = run\_\$(Process) Queue 600









- You don't have to use separate directories.
- \$(Cluster) will help distinguish runs

Universe	= vanilla
Executable	= my_job
Arguments	= -randomseed \$(Process)
Input	= my_job.input.\$(Process)
Output	<pre>= my_job.stdout.\$(Cluster).\$(Process)</pre>
Error	<pre>= my_job.stderr.\$(Cluster).\$(Process)</pre>
Log	= my_job.\$(Cluster).\$(Process).log
Queue 600	









- You ran a C program this morning
- You can also run scripting languages such as bash, python, and perl
- You can also executing programs via the command like R







- There are several different computing environments
- There is a very diverse set of computing jobs
- Matching jobs to resources is key to not wasting resources
- Not all of the available environments are open environments
- Research Computing is Complex



## Quick UNIX Refresher Before We Start



### •\$

- nano, vi, emacs, cat >, etc.
- source, module, chmod, ls









- Enough with the presentation: let's use HTCondor!
- Goal: Extend the diversity of our jobs and add some data to the mix.











- Questions? Comments?
  - Feel free to ask us questions now or later:
  - Julia Gray julia.ann.gray@gmail.com
  - Horst Severini <u>severini@ou.edu</u>
  - Pat Skubic pskubic@ou.edu
  - Exercises start here:

https://osg-htc.org/dosar/ASP2024/ASP2024\_Materials/

Presentations are also available from this URL.

