

INDIANA UNIVERSITY



The Landscape of Academic Research Computing

Kyle Gross – kagross@iu.edu
Operations Support Manager - Open Science Grid
Research Technologies - Indiana University

Some Slides Contributed by the University of Wisconsin HTCondor Team, Rob Quick, and Scot Kronenfeld



Let's jump right in...

$$\begin{split} E_{n}^{(1)} &= V_{nn} \\ E_{n}^{(2)} &= \frac{|V_{nk_2}|^2}{E_{nk_2}} \\ E_{n}^{(3)} &= \frac{|V_{nk_3}V_{k_3k_2}V_{k_2n}}{E_{nk_2}E_{nk_3}} - V_{nn} \frac{|V_{nk_3}|^2}{E_{nk_2}^2} \\ E_{n}^{(4)} &= \frac{|V_{nk_4}V_{k_4k_3}V_{k_3k_2}V_{k_2n}}{E_{nk_2}E_{nk_3}E_{nk_4}} - \frac{|V_{nk_4}|^2}{E_{nk_2}} \frac{|V_{nk_4}|^2}{E_{nk_2}E_{nk_3}E_{nk_4}} - V_{nn} \frac{|V_{nk_4}V_{k_4k_2}V_{k_2n}}{E_{nk_2}E_{nk_4}} + V_{nn}^2 \frac{|V_{nk_4}|^2}{E_{nk_4}^2} \\ &= \frac{|V_{nk_4}V_{k_4k_3}V_{k_3k_2}V_{k_2n}}{E_{nk_2}E_{nk_3}E_{nk_4}} - E_{n}^{(2)} \frac{|V_{nk_4}|^2}{E_{nk_4}^2} - 2V_{nn} \frac{|V_{nk_4}V_{k_4k_3}V_{k_3n}}{E_{nk_3}^2E_{nk_4}} + V_{nn}^2 \frac{|V_{nk_4}|^2}{E_{nk_3}^2} \\ &= \frac{|V_{nk_4}V_{k_4k_3}V_{k_3k_2}V_{k_2n}}{E_{nk_2}E_{nk_3}E_{nk_4}} - E_{n}^{(2)} \frac{|V_{nk_4}|^2}{E_{nk_4}^2} - 2V_{nn} \frac{|V_{nk_4}V_{k_4k_3}V_{k_3n}}{E_{nk_3}^2E_{nk_4}} + V_{nn}^2 \frac{|V_{nk_3}|^2}{E_{nk_3}^2} \\ &= \frac{|V_{nk_5}V_{k_5k_4}V_{k_4k_3}V_{k_3k_2}V_{k_2n}}{E_{nk_2}E_{nk_3}E_{nk_4}} - \frac{|V_{nk_5}V_{k_5k_4}V_{k_4k_3}V_{k_3n}}{E_{nk_2}E_{nk_5}} - \frac{|V_{nk_5}V_{k_5k_4}V_{k_4k_3}V_{k_3n}}{E_{nk_4}E_{nk_5}} - \frac{|V_{nk_5}V_{k_5k_4}V_{k_4k_2}V_{k_2n}}{E_{nk_2}E_{nk_5}E_{nk_5}} - V_{nn} \frac{|V_{nk_5}V_{k_5k_4}V_{k_4k_3}V_{k_3n}}{E_{nk_5}E_{nk_5}E_{nk_5}} + V_{nn} \frac{|V_{nk_5}V_{k_5k_4}V_{k_4n}}{E_{nk_5}E_{nk_$$



Who Am I?



- Operations team for Open Science Grid 10+ years
- OSG Operations Support Manager
- SWAMP Support Coordinator





Protein Docking Project at the IU School of Medicine

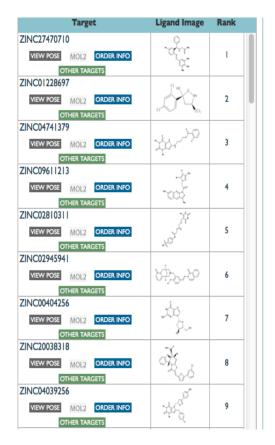


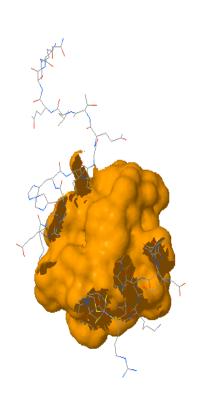
- SPLINTER Structural Protein-Ligand Interactome
- Used autodock-vina "...open-source program for drug discovery, molecular docking and virtual screening..."
- First run in 2013 docked ~3900 Proteins with 5000 Ligands for a total of ~19M docked pairs.
- Submitted via command line to Condor using Pegasus on the OSG-XSEDE submission node
- Infrastructure is set and new runs can be easily started

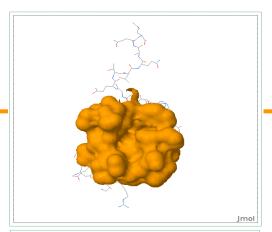


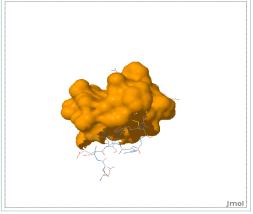


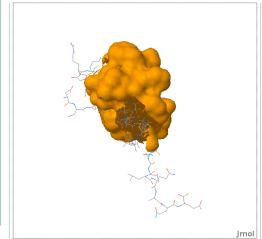
•Various rotations of Protein CBFA2T1 (Cyclin-D-related protein) (Eight twenty one protein) (Protein ETO) (Protein MTG8) (Zinc finger MYND domain-containing protein 2)









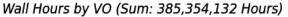


Jmol

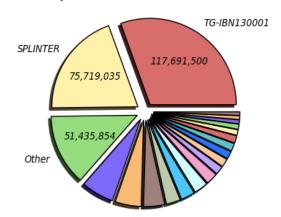


Some Numbers





1309 Days from Week 00 of 2013 to Week 31 of 2016



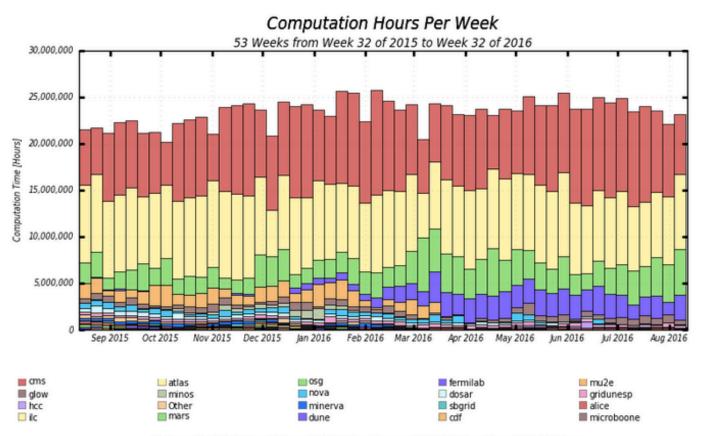
- TG-IBN130001 (117,691,501) DUKE-QGP (23,624,207) ■ CPDARKMATTERSIMULATION (10,292,491) ALGDOCK (8,520,865) SOURCECODING (5,684,844)
- DUKE-4FERMION (3,939,893) FUTURECOLLIDERS (2,684,213)
- SPLINTER (75,719,035) LIGO (18,859,313) Z2DQMC (10,049,026) RIT (5,966,501) SPHENIX (5,589,611) BIOGRAPH (3,652,186) ■ DETECTORDESIGN (2,511,013)
- Other (51,435,854) AMS (15,191,200) SNOWMASS (9,972,493) SEQ2FUN (5,961,412) NUMFPI (5,222,110) UPRRP-MR (2,786,365)
- Amazon EC2 Computing \$0.073/hour
- \$5.5M Compute Only
- Data Transfer and Storage Not Included



At what scale?



Job Activity



Maximum: 25,760,175 Hours, Minimum: 284,123 Hours, Average: 22,932,758 Hours, Current: 284,123 Hours





Follow Along at:



https://twiki.opensciencegrid.org/bin/view/ Education/ASP2016/

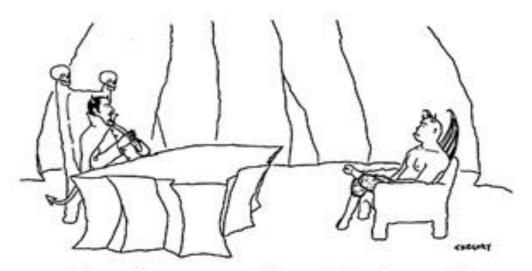




Overview of day



- Lectures alternating with exercises
 - Emphasis on lots of exercises
 - Hopefully overcome PowerPoint fatigue



"I need someone well versed in the art of torture—do you know PowerPoint?"





Some thoughts on the exercises



- 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"





INDIANA UNIVERSITY

Most important!



- Please ask questions!
 - ...during the lectures
 - ...during the exercises
 - ...during the breaks
 - ...during the meals
 - ...over dinner
 - ...via email after we depart (kagross@iu.edu or rquick@iu.edu)
- If I don't know, I'll find the right person to answer your question.



Goals for this session



- Define Local, Clustered, High Throughput Computing (HTC), High Performance Computing (HPC), and Cloud Computing (XaaS)
- Shared, Allocated, and Purchased
- What is HTCondor? And why are we using it in this Summer School?





The setup: You have a problem



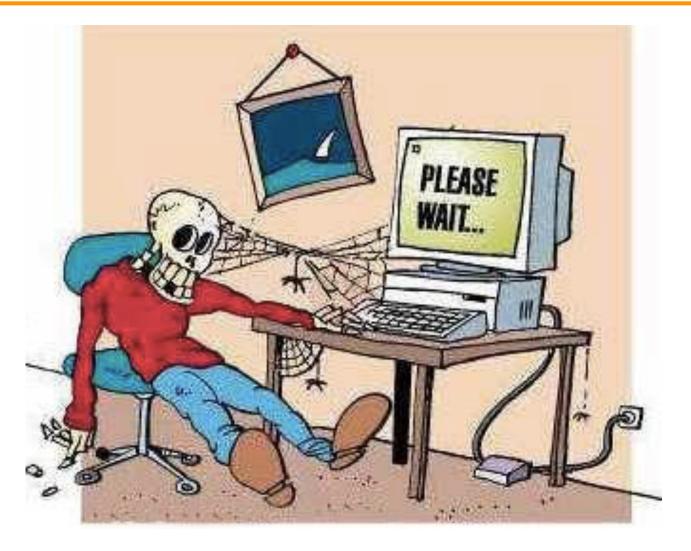
- Your science computing is complex!
 - Monte carlo, image analysis, genetic algorithm, simulation...
- It will take a year to get the results on your laptop, but the conference is in a week.
- What do you do?





Option 1: Wait a year









Option 2: Local Cluster Computing



- Easy access to additional nodes
- Local support for porting to environment (maybe)
- Often a single type of resource
- Often running at capacity





Option 3: Use a "supercomputer" aka High Performance Computing(HPC)

- "Clearly, I need the best, fastest computer to help me out"
- Maybe you do…
 - Do you have a highly parallel program?
 - i.e. individual modules must communicate
 - Do you require the fastest network/disk/memory?
- Are you willing to:
 - Port your code to a special environment?
 - Request and wait for an allocation?



INDIANA UNIVERSITY

Option 4: Use lots of commodity computers



- Instead of the fastest computer, lots of individual computers
- May not be fastest network/disk/memory, but you have a lot of them
- Job can be broken down into separate, independent pieces
 - If I give you more computers, you run more jobs
 - You care more about total quantity of results than instantaneous speed of computation
- This is high-throughput computing



Option 5: Buy (or Borrow) some Open Science Grid computing from a Cloud Provider



- Unlimited resources (if you can afford them)
- Full administrative access to OS of the resources you 'buy'
- Specialized VM images reducing effort in porting
- XaaS Business Model





These are All Valid Options



- Remember the problem: You have one month to publish results for your conference
 - Option 1: You will miss your deadline
 - Option 2: You might miss your deadline But if you're lucky you'll make it (or if you know the admin)
 - Option 3: If you have parallelized code and can get an allocation you have a good chance
 - Option 4: If you can serialize your workflow you have a good chance
 - Option 5: You can meet your deadline for a price. Though some efforts are underway to enable academic clouds





INDIANA UNIVERSITY

Computing Infrastructures



- Local Laptop/Desktop Short jobs with small data
- Local Cluster Larger jobs and larger data but subject to availability
- HPC Prime performance with parallelized code
- HTC Sustained computing over a long period for serialized
- Cloud Need deeper permission on an OS and have deeper pockets



Why focus on high-throughput computing? (HTC)



- An approach to distributed computing that focuses on long-term throughput, not instantaneous computing power
 - We don't care about operations per second
 - We care about operations per year
- Implications:
 - Focus on reliability
 - Use all available resources



Any Linux based machine can participate



Think about a race



- Assume you can run a three minute kilometer
- Does that mean you can run a 126 minute marathon?
- The challenges in sustained computation are different than achieving peak in computation speed







Why is HTC hard?



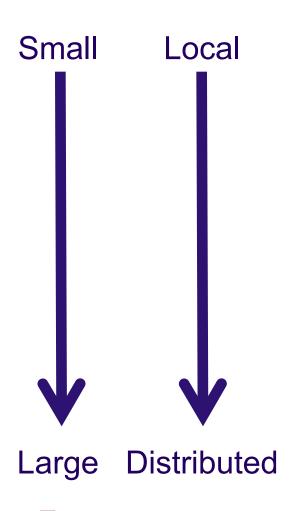
- The HTC system has to keep track of:
 - Individual tasks (a.k.a. jobs) & their inputs
 - Computers that are available
- The system has to recover from failures
 - There will be failures! Distributed computers means more chances for failures.
- You have to share computers
 - Sharing can be within an 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!)





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





Discussion



- For 5 minutes, talk to a neighbor: If you want to run one job in a local environment:
 - 1) What do you (the user) need to provide so a single job can be run?
 - 2) What does the system need to provide so your single job can be run?
 - Think of this as a set of processes: what needs happen when the job is given? A "process" could be a computer process, or just an abstract task.





What does the user provide?



- 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





What does the system provide?



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



Quick UNIX Refresher Before We Start



- \$ #This symbolizes the prompt.
- ssh <u>username@</u>training.osgconnect.net
- nano, vi, emacs, cat >, etc.
- which, rpm, ps, mkdir, cd, gcc, ls
- A variety of condor_* commands





Questions?



- Questions? Comments?
 - Feel free to ask me questions now or later:
 Kyle Gross kagross@iu.edu

Exercises start here:

https://twiki.grid.iu.edu/bin/view/Education/ RDASummerSchool/RDASchoolMaterials

Presentations are also available from this URL.

