Supercomputing in Plain English

High Throughput Computing

Henry Neeman, Director

OU Supercomputing Center for Education & Research

University of Oklahoma Information Technology

Tuesday April 19 2011

This is an experiment!

It's the nature of these kinds of videoconferences that **FAILURES ARE GUARANTEED TO HAPPEN! NO PROMISES!**

So, please bear with us. Hopefully everything will work out well enough.

If you lose your connection, you can retry the same kind of connection, or try connecting another way.

Remember, if all else fails, you always have the toll free phone bridge to fall back on.

Access Grid

If you aren't sure whether you have AG, you probably don't.

Tue Apr 19	Mosaic
Tue Apr 26	Monte Carlo
Tue May 3	Helium

Many thanks to Patrick Calhoun of OU for setting these up for us.

H.323 (Polycom etc)

From an H.323 device (e.g., Polycom, Tandberg, Lifesize, etc):

- If you **ARE** already registered with the OneNet gatekeeper:
  Dial
  2500409
- If you **AREN'T** registered with the OneNet gatekeeper (probably the case):
  1. Dial:
     164.58.250.47
  2. Bring up the virtual keypad.
     On some H.323 devices, you can bring up the virtual keypad by typing:
     #
  3. When asked for the conference ID, enter:
     0409
  4. On some H.323 devices, you indicate the end of conference ID with:
     #

Many thanks to Roger Holder and OneNet for providing this.

H.323 from Internet Explorer

From a Windows PC running Internet Explorer:

1. You **MUST** have the ability to install software on the PC (or have someone install it for you).
2. Download and install the latest Java Runtime Environment (JRE) from here:
   http://www.oracle.com/technetwork/java/javase/downloads/
   (Click on the Java Download icon, because that install package includes both the JRE and other components.)
3. Download and install this video decoder:
   http://164.58.250.47/codian video decoder.msi
4. Start Internet Explorer.
5. Copy-and-paste this URL into your IE window:
   http://164.58.250.47/
6. When that webpage loads, in the upper left, click on "Streaming."
7. In the textbox labeled Sign-in Name, type your name.
8. In the textbox labeled Conference ID, type this:
   0409
9. Click on "Stream this conference."
10. When that webpage loads, you may see, at the very top, a bar offering you options. If so, click on it and choose "Install this add-on."

H.323 from XMeeting (MacOS)

From a Mac running MacOS X:

1. Download XMeeting from
   http://xmeeting.sourceforge.net/
2. Install XMeeting as follows:
   a. Open the .dmg file.
   b. Drag XMeeting into the Applications folder.
3. Open XMeeting from Applications.
4. Skip the setup wizard.
5. In the call box, type
   164.58.250.47
6. Click the **Call** button.
7. From the Remote Control window, when prompted to join the conference, enter :
   0409#

EVO

There's a quick tutorial on the OSCER education webpage.

QuickTime Broadcaster

If you cannot connect via the Access Grid, H.323 or iLinc, then you can connect via QuickTime:

`rtsp://129.15.254.141/test_hpc09.sdp`

We recommend using QuickTime Player for this, because we've tested it successfully.

We recommend upgrading to the latest version at:

`http://www.apple.com/quicktime/`

When you run QuickTime Player, traverse the menus

File -> Open URL

Then paste in the rstp URL into the textbox, and click OK.

Many thanks to Kevin Blake of OU for setting up QuickTime Broadcaster for us.

WebEx

We have only a limited number of WebEx connections, so please avoid WebEx unless you have **NO OTHER WAY TO CONNECT**.

Instructions are available on the OSCER education webpage.

Thanks to Tim Miller of Wake Forest U.

Phone Bridge

If all else fails, you can call into our toll free phone bridge:

US: 1-800-832-0736, *6232874#

International: 303-330-0440, *6232874#

Please mute yourself and use the phone to listen.

Don't worry, we'll call out slide numbers as we go.

Please use the phone bridge **ONLY** if you cannot connect any other way: the phone bridge is charged per connection per minute, so our preference is to minimize the number of connections.

Many thanks to Amy Apon and U Arkansas for providing the previous toll free phone bridge.

Please Mute Yourself

No matter how you connect, please mute yourself, so that we cannot hear you.

At OU, we will turn off the sound on all conferencing technologies.

That way, we won't have problems with echo cancellation.

Of course, that means we cannot hear questions.

So for questions, you'll need to send some kind of text.

Questions via Text: iLinc or E-mail

Ask questions via e-mail to **sipe2011@yahoo.com**.

All questions will be read out loud and then answered out loud.

Thanks for helping!

- OSCER operations staff: Brandon George, Dave Akin, Brett Zimmerman, Josh Alexander
- Horst Severini, OSCER Associate Director for Remote & Heterogeneous Computing
- OU Research Campus staff (Patrick Calhoun, Mark McAvoy)
- Kevin Blake, OU IT (videographer)
- John Chapman, Jeff Pummill and Amy Apon, U Arkansas
- James Deaton and Roger Holder, OneNet
- Tim Miller, Wake Forest U
- Jamie Hegarty Schwettmann, i11 Industries

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Supercomputing Exercises

Want to do the "Supercomputing in Plain English" exercises?

- The first exercise is already posted at:

  http://www.oscer.ou.edu/education.php

- If you don't yet have a supercomputer account, you can get a temporary account, just for the "Supercomputing in Plain English" exercises, by sending e-mail to:

  hneeman@ou.edu

Please note that this account is for doing the **exercises only**, and will be shut down at the end of the series.

- This week's N-Body exercise will give you experience using a Condor pool.

Summer Workshops 2011

- In Summer 2011, there will be several workshops on HPC and Computational and Data Enabled Science and Engineering (CDESE) across the US.
- These will be weeklong intensives, running from Sunday evening through Saturday morning.
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Texas Advanced Computing Center
U. Texas Austin

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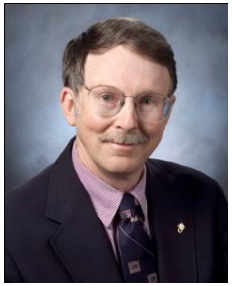

2009 Keynote:
Douglass Post
Chief Scientist
US Dept of Defense HPC Modernization Program

2010 Keynote:
Horst Simon
Deputy Director
Lawrence Berkeley National Laboratory

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2011 Keynote to be announced

FREE! Wed Oct 12 2011 @ OU

http://symposium2011.oscer.ou.edu/

Parallel Programming Workshop
FREE! Tue Oct 11 2011 @ OU
FREE! Symposium Wed Oct 12 2011 @ OU

SC11 Education Program

- At the SC11 supercomputing conference, we'll hold our annual Education Program, Sat Nov 12 – Tue Nov 15.
- You can apply to attend, either fully funded by SC11 or self-funded.
- Henry is the SC11 Education Chair.
- We'll alert everyone once the registration website opens.

Outline

- What is High Throughput Computing?
- Tightly Coupled vs Loosely Coupled
- What is Opportunistic Computing?
- Condor
- Grid Computing

What is High Throughput Computing?

High Throughput Computing

High Throughput Computing (HTC) means getting lots of work done per large time unit (for example, jobs per month).

This is different from **High *Performance* Computing** (HPC), which means getting **a particular job** done in less time (for example, calculations per second).

Throughput vs Performance

- **Throughput** is a side effect of how much time your job takes from when you first submit it until it completes.
- **Performance** is the factor that controls how much time your jobs takes from when it first starts running until it completes.
- Example:
  - You submit a very big job at 1:00am on January 1.
  - It sits in the queue for a while.
  - It starts running at 5:00pm on January 2.
  - It finishes running at 6:00pm on January 2.
  - Its performance is fast; its throughput is slow.

High Throughput on a Cluster?

Is it possible to get high throughput on a cluster?

Sure – it just has to be a cluster that no one else is trying to use!

Normally, a cluster that is shared by many users is fully loaded with jobs all the time. So your throughput depends on when you submit your jobs, and even how many jobs you submit at a time.

Depending on a variety of factors, a job you submit may wait in the queue for anywhere from seconds to days.

Tightly Coupled vs Loosely Coupled

Tightly Coupled vs Loosely Coupled

- ***Tightly coupled*** means that all of the parallel tasks have to advance forward in lockstep, so they have to communicate frequently.
- ***Loosely coupled*** means that the parallel tasks can largely or completely ignore each other (little or no communication), and they can advance at different rates.

Tightly Coupled Example

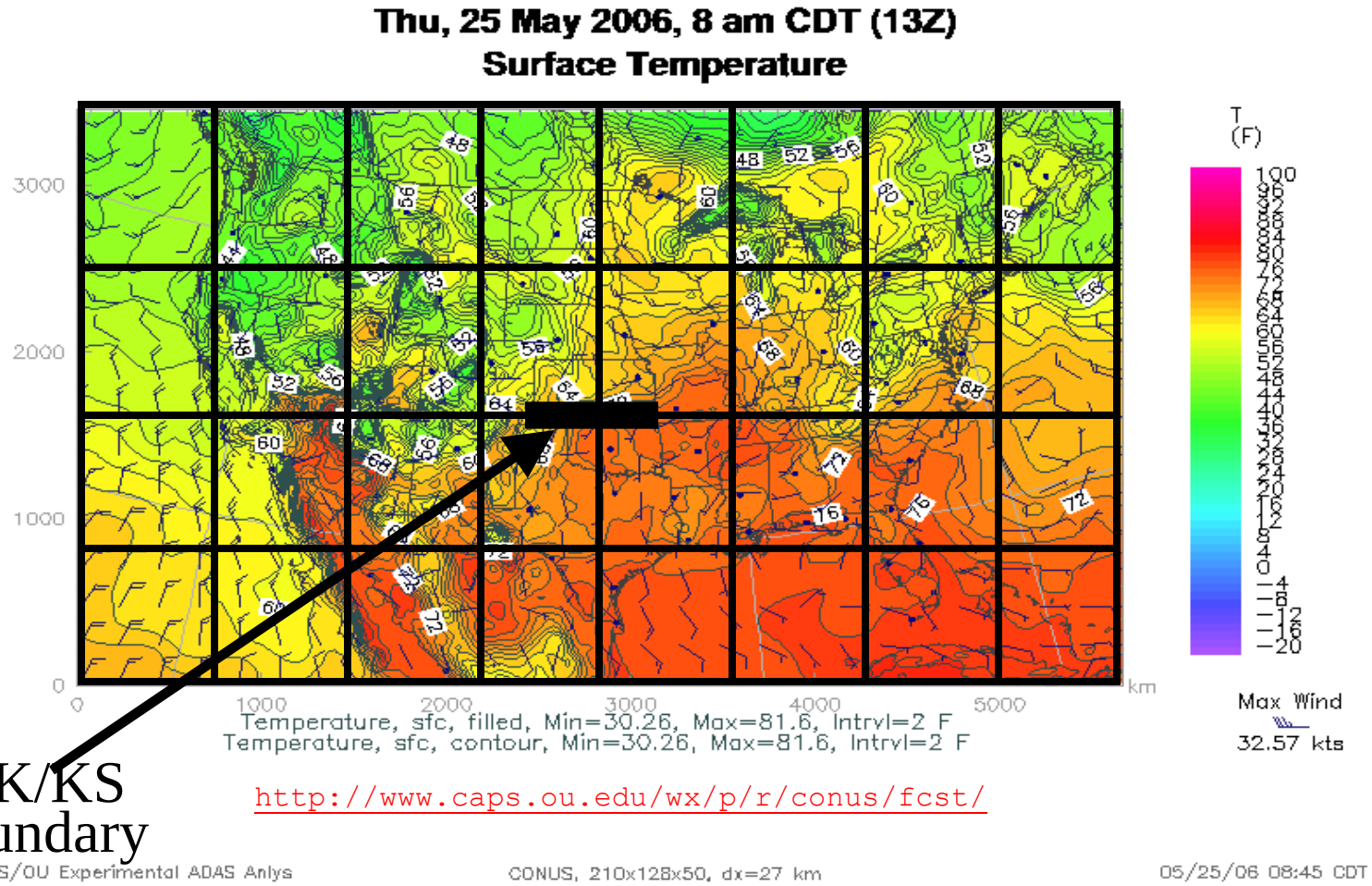
Consider weather forecasting.

You take your simulation domain – for example, the continental United States – split it up into chunks, and give each chunk to an MPI process.

But, the weather in northern Oklahoma affects the weather in southern Kansas.

So, every single timestep, the process that contains northern Oklahoma has to communicate with the process that contains southern Kansas, so that the interface between the processes has the same weather at the same time.

Tightly Coupled Example



Loosely Coupled Example

An application is known as ***embarrassingly parallel***, or ***loosely coupled***, if its parallel implementation:

1. can straightforwardly be broken up into roughly equal amounts of work per processor, **AND**
2. has minimal parallel overhead (for example, communication among processors).

We **love** embarrassingly parallel applications, because they get **near-perfect parallel speedup**, sometimes with only modest programming effort.

Monte Carlo Methods

Monte Carlo is a city in the tiny European country Monaco.

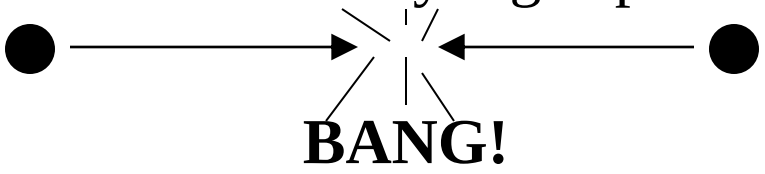
People gamble there; that is, they play games of chance, which involve randomness.

Monte Carlo methods are ways of simulating (or otherwise calculating) physical phenomena based on randomness.

Monte Carlo simulations typically are embarrassingly parallel.

Monte Carlo Methods: Example

Suppose you have some physical phenomenon. For example, consider High Energy Physics, in which we bang tiny particles together at incredibly high speeds.



We want to know, say, the average properties of this phenomenon.

There are infinitely many ways that two particles can be banged together.

So, we can't possibly simulate all of them.

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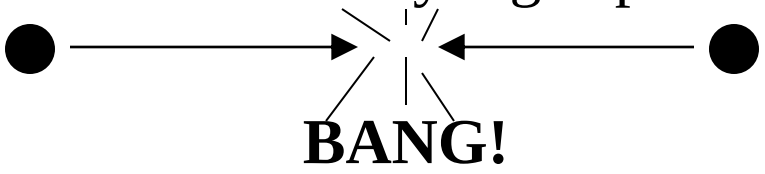

We want to know, say, the average properties of this phenomenon.

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So, we can't possibly simulate all of them.

Instead, we can **randomly choose a finite subset** of these infinitely many ways and simulate only the subset.

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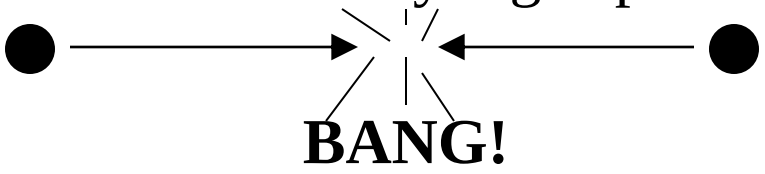


We want to know, say, the average properties of this phenomenon.

There are infinitely many ways that two particles can be banged together.

So, we can't possibly simulate all of them.

The average of this subset will be **close** to the actual average.

Monte Carlo Methods

In a Monte Carlo method, you randomly generate a large number of example cases (***realizations***) of a phenomenon, and then take the average of the properties of these realizations.

When the realizations' average **converges** (that is, doesn't change substantially if new realizations are generated), then the Monte Carlo simulation stops.

This can also be implemented by picking a high enough number of realizations to be sure, mathematically, of convergence.

MC: Embarrassingly Parallel

Monte Carlo simulations are embarrassingly parallel, because each realization is completely independent of all of the other realizations.

That is, if you're going to run a million realizations, then:

1. you can straightforwardly break up into roughly 1M / N_p chunks of realizations, one chunk for each of the N_p processes, **AND**
2. the only parallel overhead (for example, communication) comes from tracking the average properties, which doesn't have to happen very often.

Serial Monte Carlo

Suppose you have an existing serial Monte Carlo simulation:

```
PROGRAM monte_carlo
  CALL read_input(…)
  DO realization = 1, number_of_realizations
    CALL generate_random_realization(…)
    CALL calculate_properties(…)
  END DO
  CALL calculate_average(…)
END PROGRAM monte_carlo
```

How would you parallelize this?

Parallel Monte Carlo: MPI

```
PROGRAM monte_carlo_mpi
  [MPI startup]
  IF (my_rank == server_rank) THEN
    CALL read_input(…)
  END IF
  CALL MPI_Bcast(…)
  number_of_realizations_per_process = &
 &  number_of_realizations / number_of_processes
  DO realization = 1, number_of_realizations_per_process
    CALL generate_random_realization(…)
    CALL calculate_realization_properties(…)
    CALL calculate_local_running_average(...)
  END DO
  IF (my_rank == server_rank) THEN
     [collect properties]
  ELSE
     [send properties]
  END IF
  CALL calculate_global_average_from_local_averages(…)
  CALL output_overall_average(...)
  [MPI shutdown]
END PROGRAM monte_carlo_mpi
```

Parallel Monte Carlo: HTC

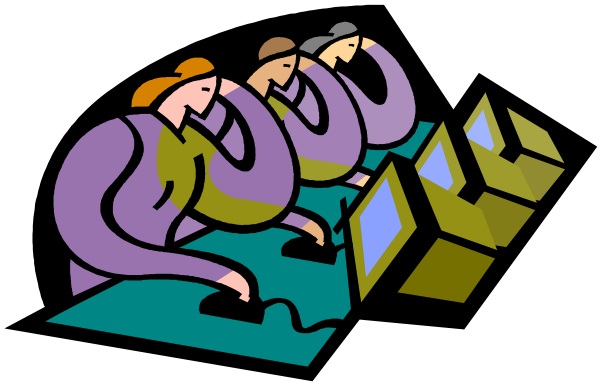
Suppose you have an existing serial Monte Carlo simulation:

```
PROGRAM monte_carlo
  CALL read_input(…)
  number_of_realizations_per_job = &
 &    number_of_realizations / number_of_jobs
  DO realization = 1, number_of_realizations_per_job
    CALL generate_random_realization(…)
    CALL calculate_properties(…)
  END DO
  CALL calculate_average_for_this_job(…)
  CALL output_average_for_this_job(…)
END PROGRAM monte_carlo
```

To parallelize this for **HTC**, simply submit **number_of_jobs** jobs, and then at the very end run a little program to calculate the overall average.

What is Opportunistic Computing?

Desktop PCs Are Idle Half the Day



Desktop PCs tend to be active during the workday.

But at night, during most of the year, they're idle. So we're only getting **half their value** (or less).

Supercomputing at Night

A particular institution – say, OU – has lots of desktop PCs that are **idle during the evening and during intersessions**.

Wouldn't it be great to put them to work on something **useful** to our institution?

That is: What if they could pretend to be a big supercomputer **at night**, when they'd **otherwise be idle anyway**?

This is sometimes known as ***opportunistic computing***:
When a desktop PC is otherwise idle, you have an opportunity to do number crunching on it.

Supercomputing at Night Example

SETI – the Search for Extra-Terrestrial Intelligence – is looking for evidence of green bug-eyed monsters on other planets, by mining radio telescope data.

SETI@home runs number crunching software as a screensaver on idle PCs around the world (2+ million PCs in 252 countries):

http://setiathome.berkeley.edu/

There are **many similar projects**:

- folding@home (protein folding)
- climateprediction.net
- Einstein@Home (Laser Interferometer Gravitational wave Observatory)
- Cosmology@home
- …

BOINC

The projects listed on the previous page use a software package named BOINC (**B**erkeley **O**pen **I**nfrastructure for **N**etwork **C**omputing), developed at the University of California, Berkeley:

http://boinc.berkeley.edu/

To use BOINC, you have to insert calls to various BOINC routines into your code. It looks a bit similar to MPI:

```
int main ()
{ /* main */
    …
    boinc_init();
    …
    boinc_finish(…);
} /* main */
```

Condor

Condor is Like BOINC

- Condor **steals computing time** on existing desktop PCs **when they're idle**.
- Condor **runs in background** when no one is sitting at the desk.
- Condor allows an institution to get **much more value** out of the hardware that's **already purchased**, because there's little or no idle time on that hardware – all of the idle time is used for number crunching.

Condor is Different from BOINC

- To use Condor, **you don't need to rewrite your software** to add calls to special routines; in BOINC, you do.
- Condor **works great under Unix/Linux**, but less well under Windows or MacOS (more on this presently); BOINC works well under all of them.
- It's **non-trivial to install Condor** on your own personal desktop PC; it's straightforward to install a BOINC application such as SETI@home.

Useful Features of Condor

- **Opportunistic** computing: Condor steals time on existing desktop PCs when they're otherwise not in use.
- Condor **doesn't require any changes to the software**.
- Condor can **automatically checkpoint** a running job: Every so often, Condor saves to disk the state of the job (the values of all the job's variables, plus where the job is in the program).
- Therefore, Condor can **preempt** running jobs if more important jobs come along, or if someone sits down at the desktop PC.
- Likewise, Condor can **migrate** running jobs to other PCs, if someone sits at the PC or if the PC crashes.
- And, Condor can do all of its **I/O over the network**, so that the job on the desktop PC doesn't consume the desktop PCs local disk.

Condor Pool @ OU

OU IT has deployed a large Condor pool (795 desktop PCs in dozens of labs around campus).

OU's Condor pool provides a huge amount of computing power – more than OSCER's big cluster:

- if OU were a state, we'd be the 17th largest state in the US;
- if OU were a country, we'd be the 10th largest country in the world.

The hardware and software cost is zero, and the labor cost is modest.

Also, we've been seeing empirically that lab PCs are available for Condor jobs about 80% of the time.

Condor Limitations

- The Unix/Linux version has **more features** than Windows or MacOS, which are referred to as "clipped."
- Your code **shouldn't be parallel** to do opportunistic computing (MPI requires a fixed set of resources throughout the entire run), and it shouldn't try to do any funky communication (for example, opening sockets).
- For a Red Hat Linux Condor pool, you have to be able to **compile your code** with gcc, g++, g77 or NAG f95 (which is a Fortran90-to-C translator that then calls gcc).
- Also, depending on the PCs that have Condor on them, you may have limitations on, for example, how big your jobs' RAM footprint can be.

Running a Condor Job

Running a job on Condor pool is a lot like running a job on a cluster:

1. You compile your code using the compilers appropriate for that resource.
2. You submit a batch script to the Condor system, which decides when and where your job runs, magically and invisibly.

Sample Condor Batch Script

```
Universe     = standard
Executable   = /home/hneeman/NBody/nbody_compiled_for_condor
Notification = Error
Notify_User  = hneeman@ou.edu
Arguments    = 1000 100
Input        = /home/hneeman/NBody/nbody_input.txt
Output       = nbody_$(Cluster)_$(Process)_out.txt
Error        = nbody_$(Cluster)_$(Process)_err.txt
Log          = nbody_$(Cluster)_$(Process)_log.txt
InitialDir   = /home/hneeman/NBody/Run001
Queue
```

The batch submission command is **`condor_submit`**, used like so:

`condor_submit  nbody.condor`

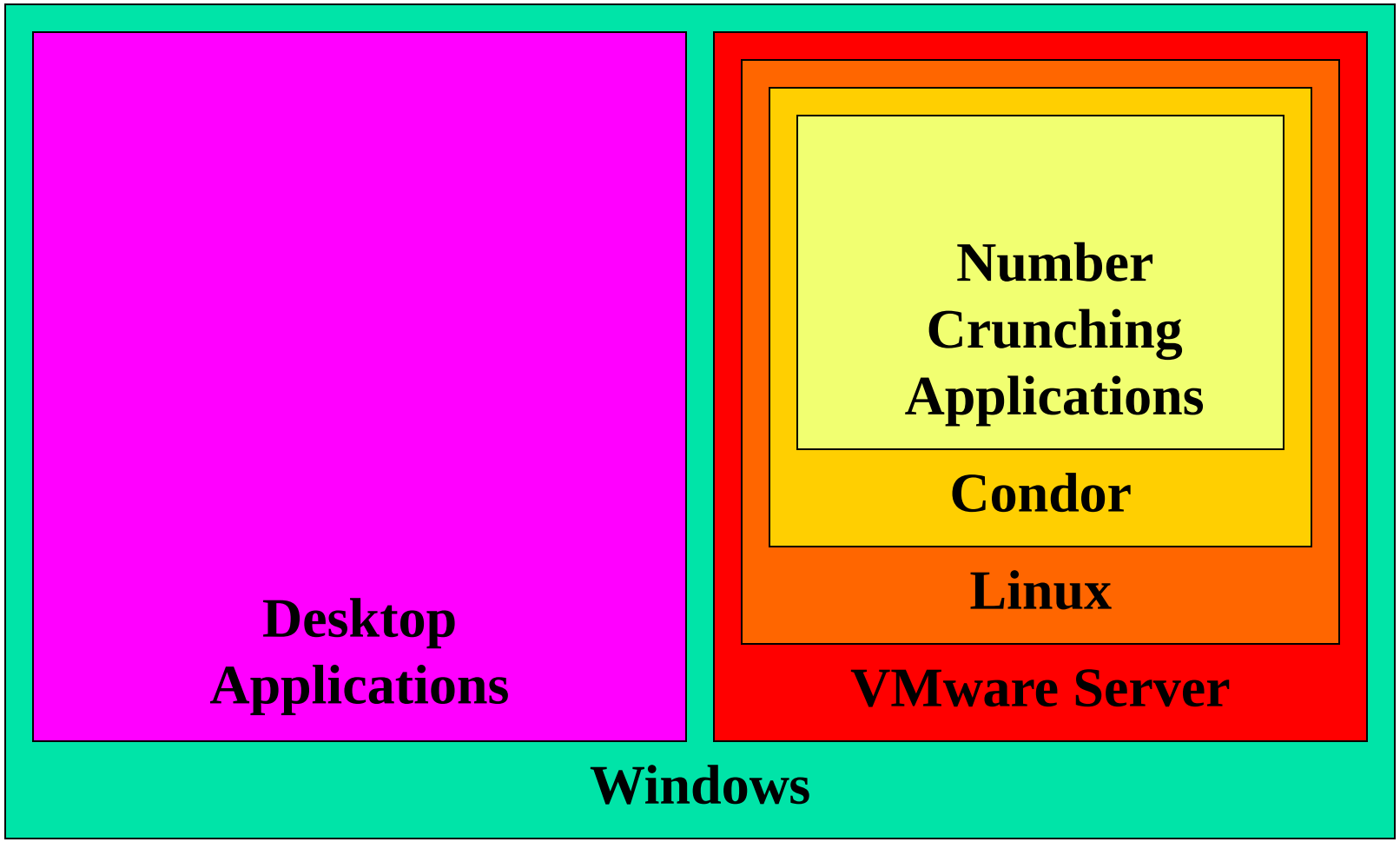
Linux Condor on Windows PCs?

If OU's Condor pool uses Linux, how can it be installed in OU IT PC labs? Don't those run Windows?

Yes.

Our solution is to run Linux inside Windows, using a piece of software named Vmware ("virtual machine"), but there are other software packages that can be used (for example, VirtualBox).

Condor inside Linux inside Windows



Advantages of Linux inside Windows

- Condor is full featured rather than clipped.
- Desktop users have a full Windows experience, without even being aware that VMware exists.
- A little kludge helps Condor watch the keyboard, mouse and CPU level of Windows, so that Condor jobs don't run when the PC is otherwise in use.

Grid Computing

What is Grid Computing?

The term ***grid computing*** is poorly defined, but the best definition I've seen so far is:

"a distributed, heterogeneous operating system."

A ***grid*** can consist of:

- compute resources;
- storage resources;
- networks;
- data collections;
- shared instruments;
- sensor networks;
- and so much more ....

Grid Computing is Like and Unlike ...

IBM's website has a very good description of grid computing:

- ***"Like the Web**, grid computing keeps complexity hidden: multiple users enjoy a single, unified experience.*
- ***"Unlike the Web**, which mainly enables communication, grid computing enables full collaboration toward common ... goals.*
- ***"Like peer-to-peer**, grid computing allows users to share files.*
- ***"Unlike peer-to-peer**, grid computing allows many-to-many sharing – not only files but other resources as well.*
- ***"Like clusters and distributed computing**, grids bring computing resources together.*
- ***"Unlike clusters and distributed computing**, which need physical proximity and operating homogeneity, grids can be geographically distributed and heterogeneous.*
- ***"Like virtualization technologies**, grid computing enables the virtualization of IT resources.*
- ***"Unlike virtualization technologies**, which virtualize a single system, grid computing enables the virtualization of vast and disparate IT resources."*

http://www.thocp.net/hardware/grid_computers.htm

Condor is Grid Computing

Condor creates a grid out of disparate desktop PCs.

(Actually, they don't have to be desktop PCs; they don't even have to be PCs. You can use Condor to schedule a cluster, or even on a big iron supercomputer.)

From a user's perspective, all of the PCs are essentially invisible; the user just knows how to submit a job, and everything happens magically and invisibly, and at some point the job is done and a result appears.

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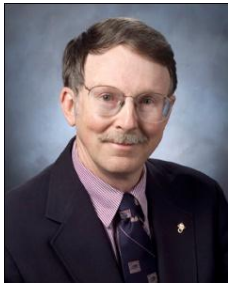

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