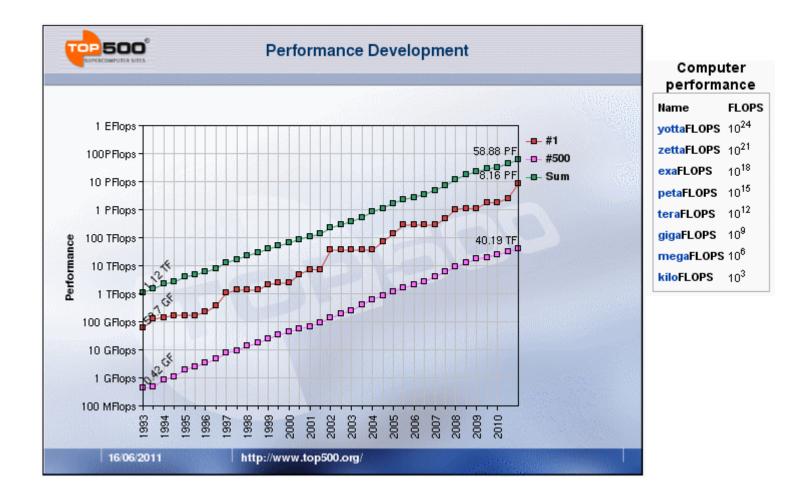
Taxes, Death, and Parallelism are Inevitable



Introduction to Parallel Computing

- <u>Abstract</u>
- Overview
 - What is Parallel Computing?
 - <u>Why Use Parallel Computing?</u>
- <u>Concepts and Terminology</u>
 - von Neumann Computer Architecture
 - Flynn's Classical Taxonomy
 - <u>Some General Parallel Terminology</u>

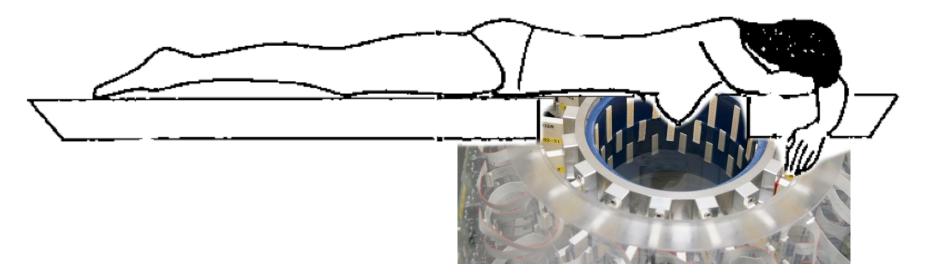
The FUTURE: *The race is already on for Exascale Computing!*



<u>Ultrasound Computer Tomography (USCT)</u> http://www.interactive-grid.eu

- New method for medical imaging
 - Focus: Breast cancer diagnosis

USCT setup



USCT Algorithm

- Characteristicts:
 - Input: 20 GB (full set)
 - Computing time depends
 - on output size / resolution
 - amount of input data

35MB	20GB	20GB	Data
4096²	128²x100	4096²x3410	Voxels
1 Hour	1.5 Months	150 Years	Time

- Matlab
 - Strategic development platform (95% sourcecode)

USCT Algorithm

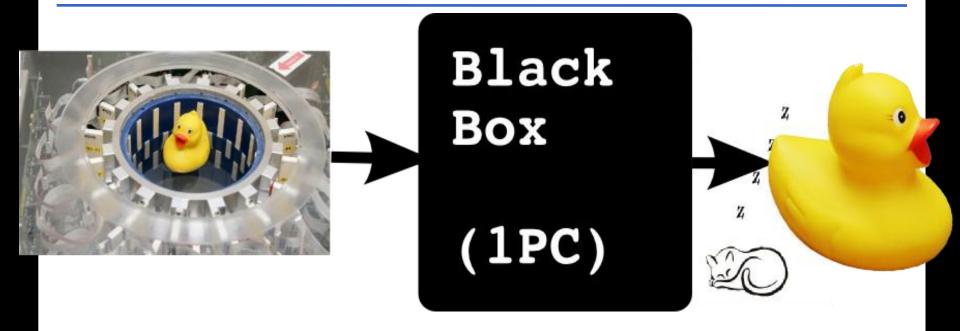
- Characteristicts:
 - Input: 20 GB (full set)
 - Computing time depends
 - on output size / resolution
 - amount of input data

35MB	20GB	20GB	Data
4096²	128²x100	4096²x3410	Voxels
1 Hour	1.5 Months	150 Years	Time



USCT reconstruction := "Black Box"





- Computation takes long (days, weeks, years)
- Grid in order to speed up

int.eu.grid http://interactive-grid.eu M

Marcus.Hardt@iwr.fzk.de





Idea: Computer power <=> Electrical power

From Electrical power grid => computational grid

- Across organisationsal domains / countries
- Transparent access to
 - Computing
 - Data

int.eu.grid

- Network
- Large scale installations



Grid middleware

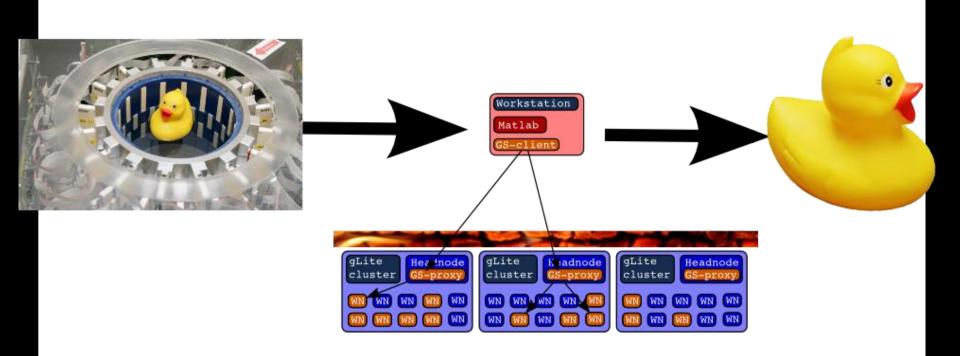


- Middleware
 - := Layer between application and operating system
- gLite: <u>one</u> grid middleware
 - Development driven by CERN
 - Tools for data+computing of new accelerator
 - 10 TB/year * 20 years, random access
- Paradigm: Send job to where the data is
- Job: Self contained application



Putting things together



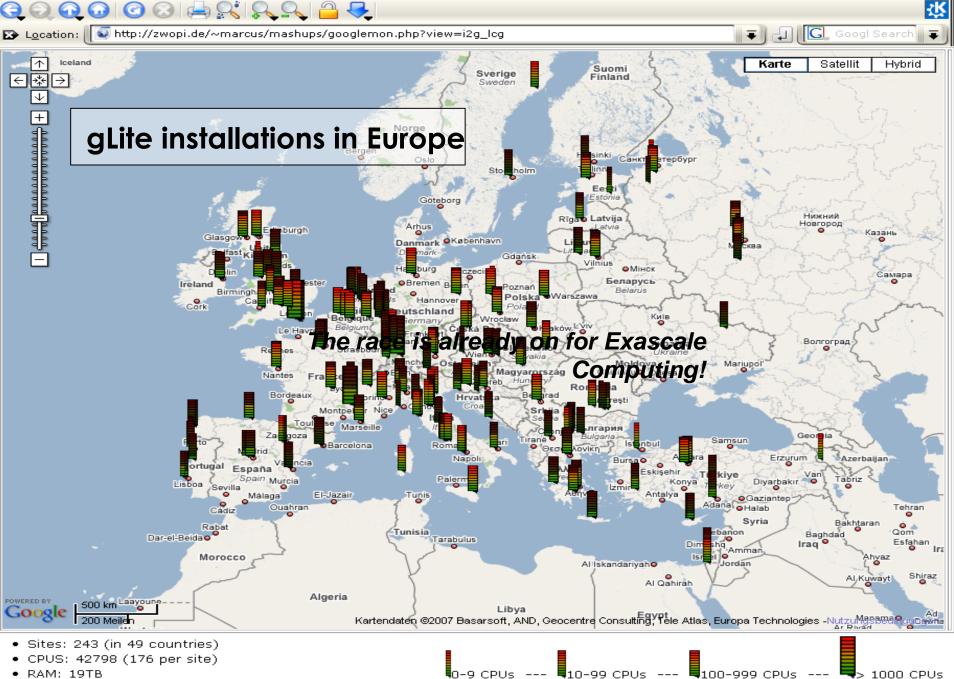


int.eu.grid

http://interactive-grid.eu

Marcus.Hardt@iwr.fzk.de



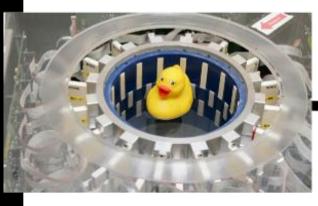


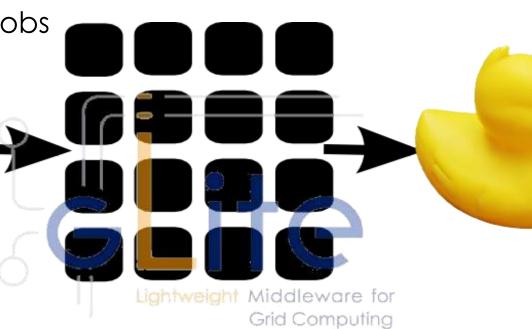
- RAM: 19TB
- RAM/CPU: 468MB
- DISK [Tot / Avail]: [8042TB / 5408TB] ([33892GB / 22792GB] per site)

Using gLite



- Initial approach to parallel execution:
 - Partitioning of data
 - Many parallel jobs

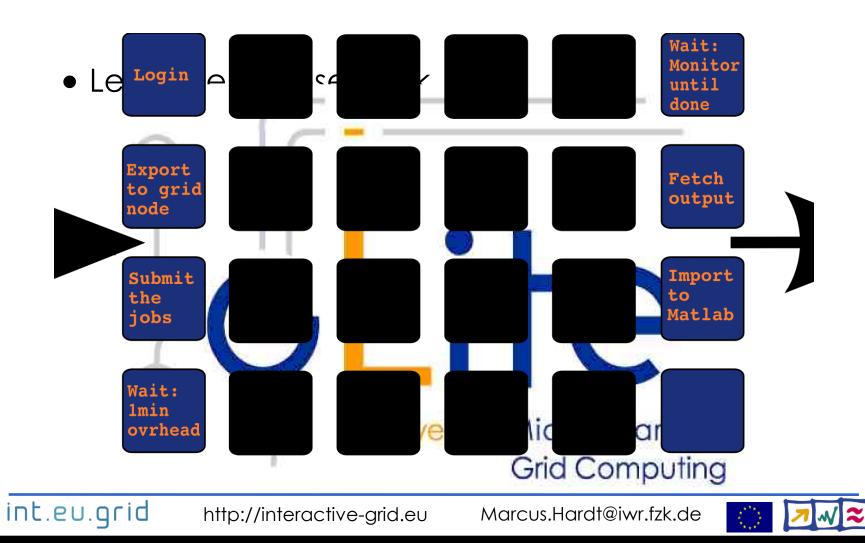


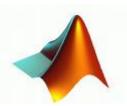




Using gLite







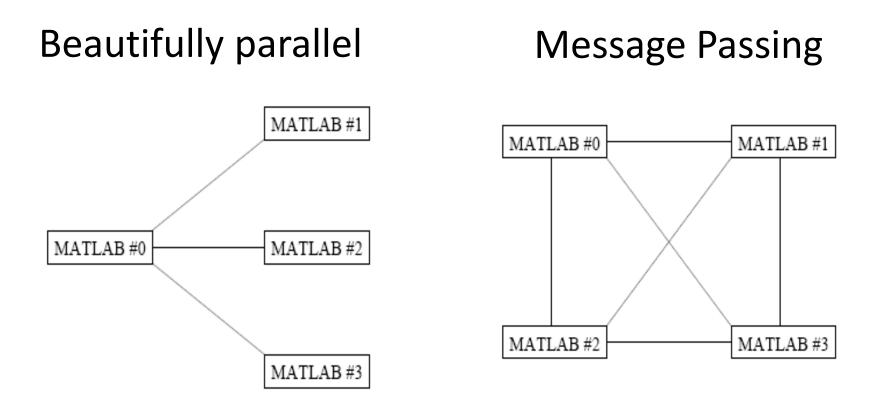
Parallel Matlab

Elias Houstis

About MATLAB

- MATLAB (with Simulink) programming language for science and engineering
- Over 1 milion users, over 3500 universities and colleges
- Engineering in industry + biotech, medical, financial
- Toolboxes for different fields
 - Engineering, Bioinformatics, Economics etc.
- Parallel computing support
 - Job execution on multicore/cluster systems
 - MPI support

Matlab in Parallel

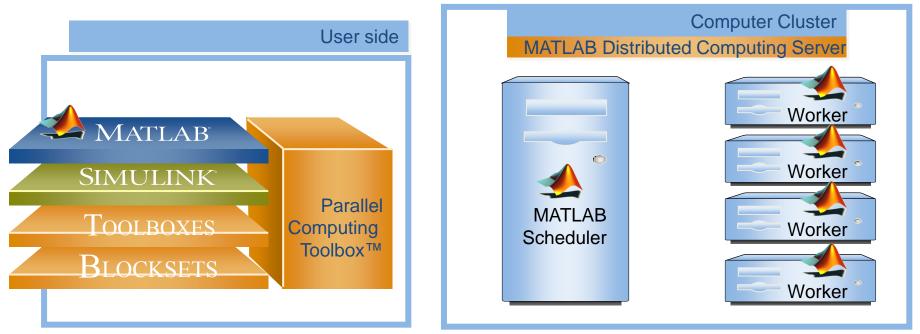


e.g., Multi, paralize, Plab, ParMatlab

e.g., MultiMatlab, CMTM, DPToolbox,MatlabMPI, pMatlab

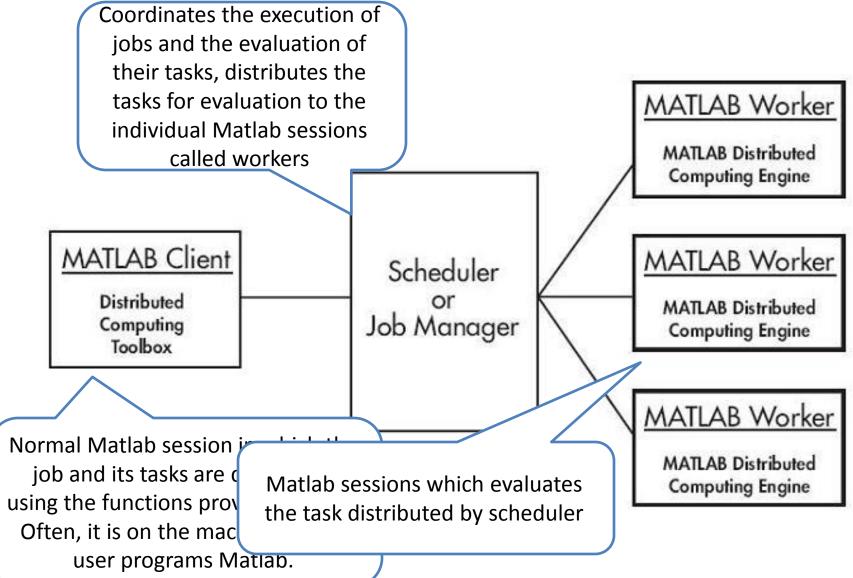
www-math.mit.edu/~edelman/homepage/papers/pmatlab.pdf

Parallel Computing With MATLAB



• Support for third party schedulers

Architecture of PCT



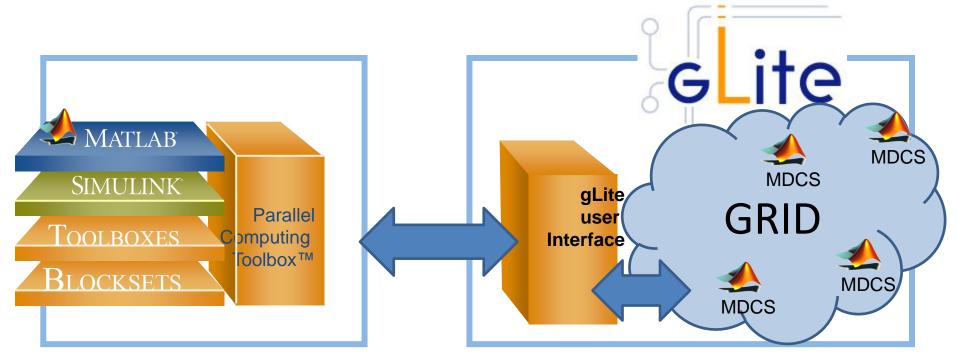
Distributed Computing Server (DCS)



- Parallel Computing Toolbox
 - Only four local workers on a multicore or multiprocessor computer
- PCT + DCS -> Cluster-based applications
- Coordinate and execute independent MATLAB operations simultaneously on a cluster of computers

MATLAB gLite integration

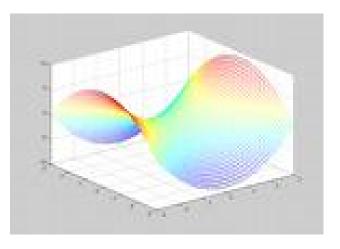
• API for generic scheduler



Developed by EGEE&Mathworks

Parallel Computing Toolbox (PCT)

- Data and task parallelism using
 - Parallel-for loops
 - Distributed arrays
 - Parallel numerical algorithms
 - Message Passing functions



• Easy transition between serial and parallel

http://www.mathworks.com/products/parallel-computing/description1.html

Mathworks – Parallel Computing toolbox

- The toolbox provides eight workers (MATLAB computational engines) to execute applications locally on a multicore desktop
- Parallel for-loops (parfor) for running taskparallel algorithms on multiple processors
- Computer cluster and grid support (with MATLAB Distributed Computing Server)

Vectorization

>> clear all;

tic;

for i=1:50000

a(i) = sin(i);

end

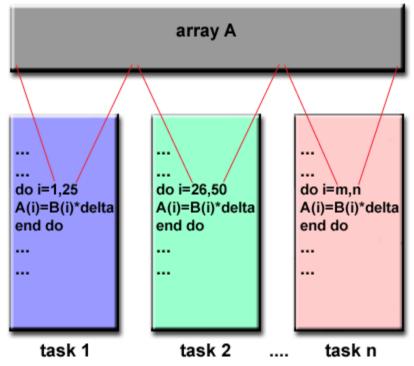
toc

Elapsed time is
3.211070 seconds.

>> clear all; tic; i = [1:50000]; a = sin(i);toc Elapsed time is 0.016062 seconds. >> speedup=3.211070 /0.016062 =199.9172

Data Parallel

- The data parallel model demonstrates the following characteristics:
 - Most of the parallel work performs operations on a data set, organized into a common structure, such as an array
 - A set of tasks works collectively on the same data structure, with each task working on a different partition
 - Tasks perform the same operation on their partition



 On shared memory architectures, all tasks may have access to the data structure through global memory. On distributed memory architectures the data structure is split up and resides as "chunks" in the local memory of each task.

parfor - Parallel for loop

parfor - Parallel for loop

Syntax

```
parfor loopvar = initval:endval; statements; end
parfor (loopvar = initval:endval, M); statements; end
```

Description

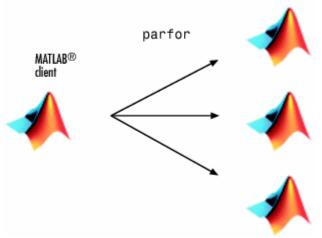
parfor loopvar = initval:endval; statements; end executes a series of MATLAB commands denoted here as statements for values of loopvar between initval and endval, inclusive, which specify a vector of increasing integer values. Unlike a traditional for-loop, there is no guarantee of the order in which the loop iterations are executed.

The general format of a parfor statement is:

```
parfor Loopvar = initval:endval
```

<statements>

end



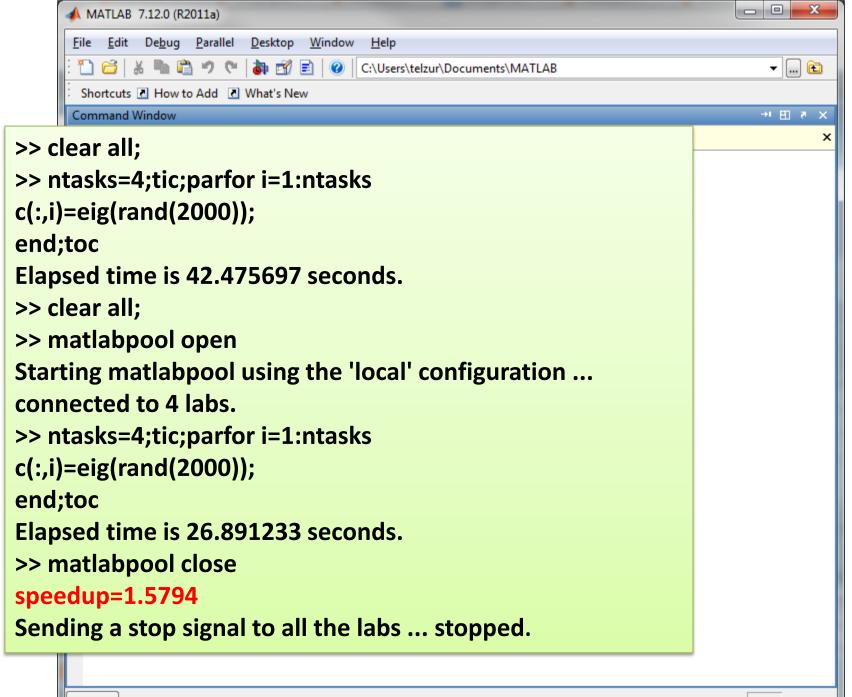
parfor – Perform three large eigenvalue computations using four computers or cores

- >> clear all;
- >> ntasks=4;tic;for
 i=1:ntasks
- c(:,i)=eig(rand(100
 0));

end;toc

Elapsed time is 8.575804 seconds.

- >> clear all;
- matlabpool open
- ntasks=4;tic;parfor i=1:ntasks
- c(:,i)=eig(rand(1000));
- end;toc
- Starting matlabpool using the 'local' configuration ... connected to 4 labs.
- Elapsed time is 5.198244 seconds.
- Speedup= 1.6522



📣 <u>S</u>tart

Parallel mode-I: matlabpool

- Open or close a pool of MATLAB sessions for parallel computation
- Syntax:

MATLABPOOL MATLABPOOL OPEN MATLABPOOL OPEN <poolsize> MATLABPOOL CLOSE MATLABPOOL CLOSE FORCE

• Work on local client PC

.....

• Without open matlabpool, parallel code will still run but runs sequentially

- %% Parameter Sweep of ODEs
- % This is a parameter sweep study of a 2nd order ODE system.
- %
- % \$m\ddot{x} + b\dot{x} + kx = 0\$
- %
- % We solve the ODE for a time span of 0 to 25 seconds, with initial
- % conditions \$x(0) = 0\$ and \$\dot{x}(0) = 1\$. We sweep the parameters \$b\$
- % and \$k\$ and record the peak values of \$x\$ for each condition. At the end,
- % we plot a surface of the results.

- Computing in serial...
- Elapsed time is 27.64 seconds.
- Computing in parallel...
- Starting matlabpool using the 'local' configuration ... connected to 4 labs.
- Elapsed time is 12.39 seconds.
- Sending a stop signal to all the labs ... stopped.
- Speed up (time serial / time parallel): 2.23

%% Parameter Sweep (Parallel) % Next, we convert the |for| loop to a |parfor| loop and start a pool or % MATLAB workers.

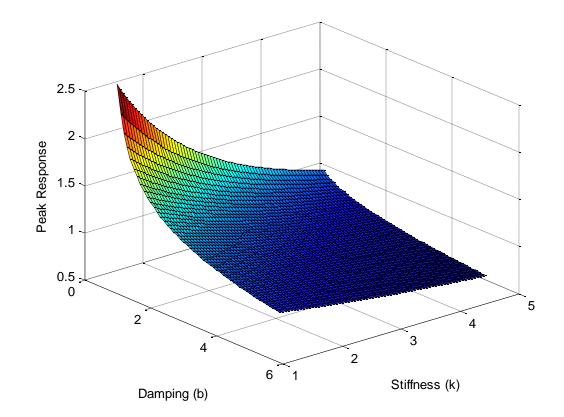
disp('Computing in parallel...');drawnow;

matlabpool open

tic; parfor idx = 1:numel(kGrid) % Solve ODE [T,Y] = ode45(@(t,y) odesystem(t, y, m, bGrid(idx), kGrid(idx)), ... [0, 25], ... % simulate for 25 seconds [0, 1]); % initial conditions

% Determine peak value peakVals(idx) = max(Y(:,1)); end t2 = toc; fprintf('Elapsed time is %0.2f seconds.\n', t2);

% Close MATLAB Pool matlabpool close



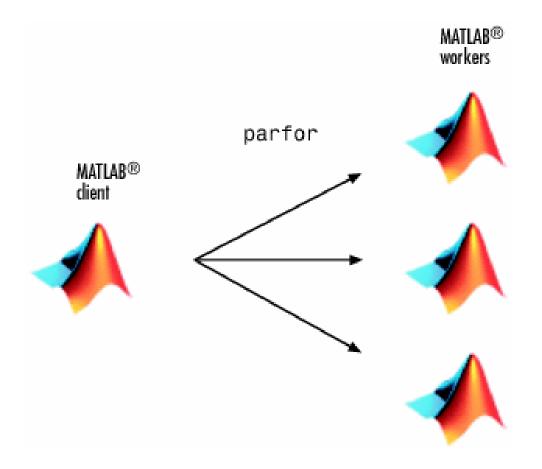
Task Parallel applications

- parallel problems by organizing them into independent *tasks* (units of work)
 - parallelize Monte Carlo simulations
- Parallel for-Loops (parfor)

```
parfor (i = 1 : n)
% do something with i
end
```

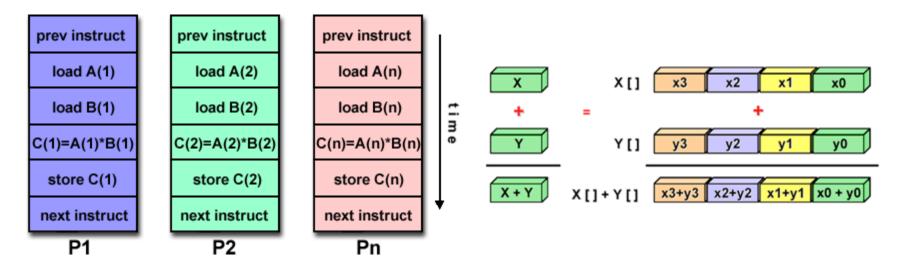
- Mix task parallel and serial code in the same function
- Run loops on a pool of MATLAB resources
- Iterations must be order-independent

Iterations run in parallel in the MATLAB pool (local workers)





- A type of parallel computer
- All processing units execute the same instruction at any given clock cycle
- Each processing unit can operate on a different data element
- Two varieties: Processor Arrays and Vector Pipelines
- Most modern computers, particularly those with graphics processor units (GPUs) employ SIMD instructions and execution units.



Data Parallel applications

 Single Program Multiple Data (spmd) spmd (n) <statements>

end

For example, create a random matrix on four labs:

matlabpool open spmd (2) R = rand(4,4); end matlabpool close create different sized arrays depending on labindex:

> matlabpool open spmd (2) if labindex==1 R = rand(4,4); else R = rand(2,2); end end matlabpool close

Demo

```
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    >> matlabpool
Current Folder Command History
   Starting matlabpool using the 'local' configuration ... connected to 2 labs.
   >> spmd(2)
   R=rand(4,4);
    end
    >> R(:)
    ans =
        [4x4 double]
        [4x4 double]
    >> R{1}
    ans =
        0.9173 0.4612 0.2155 0.4621
        0.6839 0.1562 0.4978 0.9846
               0.4626 0.2904 0.9587
        0.8661
        0.4809
                0.8009
                          0.9071
                                     0.5795
    >> R{2}
    ans =
        0.2951
                0.7010
                          0.9143 0.7375
                                                                                   E
                0.3821 0.2740 0.5407
        0.0990
        0.3277
                 0.9602 0.6484
                                     0.6348
                0.7780
                          0.2781 0.0948
        0.6902
 f_{x} >>
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```

Demo

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Shortcuts 🔄 How to Add 💽 What's New
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Current Folder Command History
     >> spmd (2)
        if labindex==1
           R = rand(4, 4);
      else
           R = rand(2,2);
        end
     end
     >> R(:)
     ans =
        [4x4 double]
         [2x2 double]
     >> R\{1\}
     ans =
         0.5324 0.9413 0.6332 0.8776
         0.1019 0.2410 0.0554 0.9100
         0.5341 0.3698 0.7112 0.5573
         0.2081 0.0299 0.0759 0.1785
     >> R\{2\}
     ans =
                                                                                       E
         0.3802 0.7370
         0.1215 0.3625
  f_{x} >>
                                                                                   OVR
📣 Start
```

Distributed arrays and operations (matlabpool mode) codistributor()

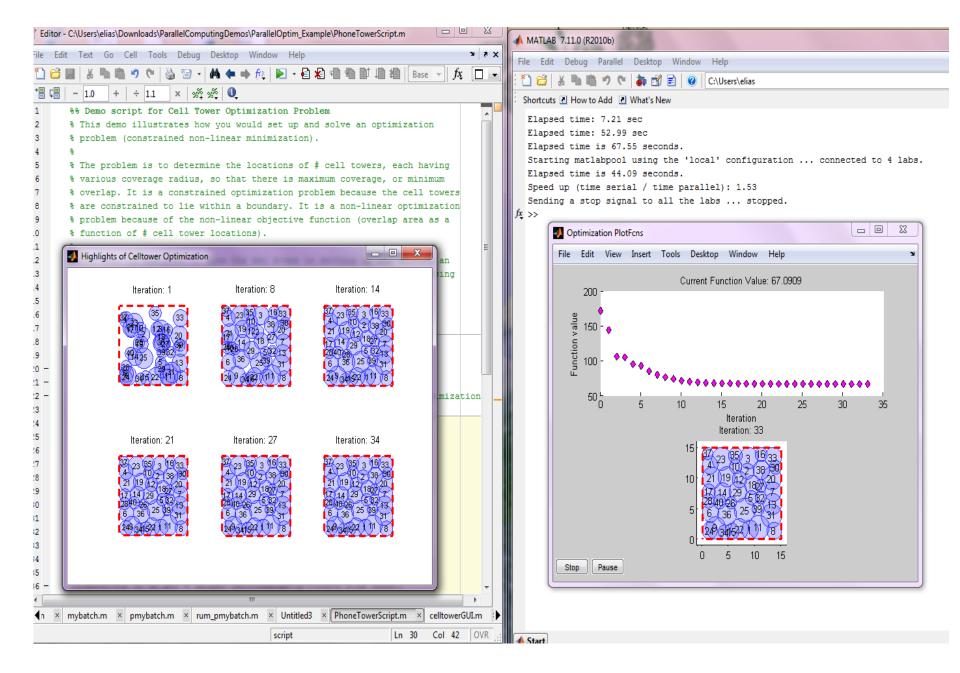
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Current Folder Command History
     >> matlabpool
     Starting matlabpool using the 'local' configuration ... connected to 2 labs.
     >> spmd
     a=rand(500,500);
     end
     >> a
     a =
        Lab 1: class = double, size = [500
                                                5001
        Lab 2: class = double, size = [500
                                                5001
     >> spmd
     b=rand(800,800, codistributor())
     end
     Lab 1:
       This lab stores b(:,1:400).
                LocalPart: [800x400 double]
           Codistributor: [1x1 codistributor1d]
     Lab 2:
       This lab stores b(:,401:800).
                LocalPart: [800x400 double]
           Codistributor: [1x1 codistributor1d]
  f_{\underline{x}} >>
                                                                                              OVR
  Start
```

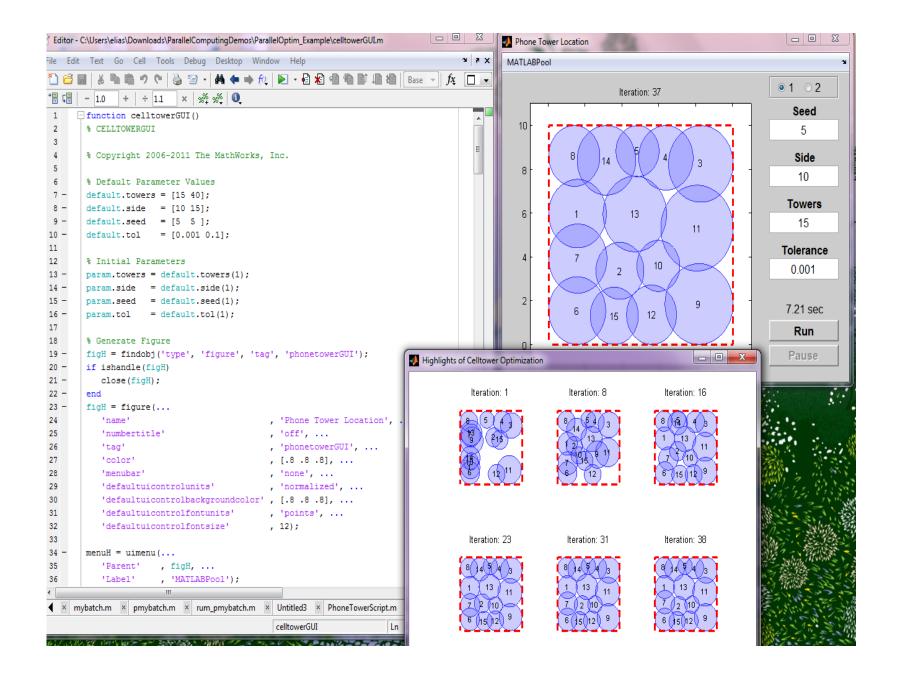
```
- 23
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Shortcuts How to Add What's New
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Current Folder Command History
    >> spmd
    c=b'
    end
    Lab 1:
      This lab stores c(1:400,:).
              LocalPart: [400x800 double]
         Codistributor: [1x1 codistributor1d]
    Lab 2:
      This lab stores c(401:800,:).
              LocalPart: [400x800 double]
          Codistributor: [1x1 codistributor1d]
    >> c
    c = \langle page truncated: showing [1:32, 1:32] of 800-by-800 \rangle
      Columns 1 through 8
        0.2272 0.5522 0.9372 0.8589 0.8029 0.4314 0.2116
                                                                              0.5545
        0.0658 0.7758 0.0901 0.3114 0.0746 0.5446 0.5271
                                                                              0.7786
        0.0811
               0.2854 0.2272 0.1215
                                                          0.9134
                                                                              0.7143
                                                0.3944
                                                                    0.6010
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Shortcuts 🔄 How to Add 🔃 What's New
    >>
Current Folder Command History
                                                                                             a.
    >> spmd
    d=c*b
    end
    Lab 1:
      This lab stores d(:,1:400).
               LocalPart: [800x400 double]
          Codistributor: [1x1 codistributor1d]
    Lab 2:
       This lab stores d(:,401:800).
               LocalPart: [800x400 double]
           Codistributor: [1x1 codistributor1d]
    >> d
    d = <page truncated: showing [1:32, 1:32] of 800-by-800>
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codistributed()

```
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Command History
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     >> spmd
     a=rand(500,500);
     b=codistributed(a)
     end
     Lab 1:
Current Folder
       This lab stores b(:,1:250).
                LocalPart: [500x250 double]
            Codistributor: [1x1 codistributor1d]
     Lab 2:
       This lab stores b(:,251:500).
                LocalPart: [500x250 double]
            Codistributor: [1x1 codistributor1d]
     >>
     >>
  f_{x} >>
      ₹.
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                                                                                           OVR
```





• Parallel mode on a MATLAB Pool (1)

matlabpool Open or close pool of MATLAB sessions for parallel computation
parfor Execute code loop in parallel
spmd Execute code in parallel on MATLAB pool

Interactive Functions

help Help for toolbox functions in Command Window pmode Interactive Parallel Command Window

Parallel mode on a MATLAB Pool (2)

batch Run MATLAB script as batch job

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*= C= - 1.0 + ÷ 1.1 × % % 0	Shortcuts 🗷 How to Add 🗷 What's New
This file uses Cell Mode. For information, see the rapid code.	>> [time1,time2,time3] = newSPMDVersion
1 [function [time1,time2,time3] = new	Starting matlabpool using the 'local' configuration connected to 4 labs.
2 - n=10^8; step = 1/n;	Sending a stop signal to all the labs stopped.
3 % Serial Version	timel =
4 - tic; s=0;	
<pre>5 = G for i=1:n-1 6 = x=(i-0.5)*step; s=s+4./(1+x^2);</pre>	2.7594
7 end	
8 - timel = toc;	
9	time2 =
10 % Parallel Version	1.0683
11 - matlabpool open 4	
12 - tic;	
<pre>13 - spmd 14 - slocal = myTestSum(n, step);</pre>	time3 =
15 end	
16 - time2 = toc;	0.9888
17 % Parallel Version parfor	fx >>
<pre>18 - nstep=10^8; step = 1/nstep;</pre>	
19 - tic;s=0;	
20 - parfor i=1:nstep-1	
21 - x=(1-0.5)*step; 22 - s=s+4./(1+x^2);	
23 - end	
24 - time3 = toc;matlabpool close	
25 end	
26 %%	
27 Efunction slocal = myTestSum(n, step	
28 - nlo = (n * (labindex - 1)) / n	
<pre>29 - nhi = (n * labindex) / numlabs; 30 - slocal = 0;</pre>	
31 - for i=nlo:nhi	
32 - x=(1-0.5)*step; slocal=slocal+4./(
33 - end	
34 - end	

function [time1,time2,time3] = newSPMDVersion $n=10^8$; step = 1/n; % Serial Version tic: s=0; for i=1:n-1x=(i-0.5)*step; s=s+4./(1+x^2); end time1 = toc;% Parallel Version matlabpool open 4 tic; spmd slocal = myTestSum(n,step); end time2 = toc;

% Parallel Version parfor nstep=10^8; step = 1/nstep; tic;s=0; parfor i=1:nstep-1 x=(i-0.5)*step; s=s+4./(1+x^2); end time3 = toc;matlabpool close end

```
function slocal = myTestSum(n,step)
nlo = ( n * ( labindex - 1 ) ) / numlabs + 1;
nhi = ( n * labindex) / numlabs;
slocal = 0;
for i=nlo:nhi
x=(i-0.5)*step; slocal=slocal+4./(1+x^2);
end
end
```

Profiling (profile on...profile viewer)

- profile on
- computations

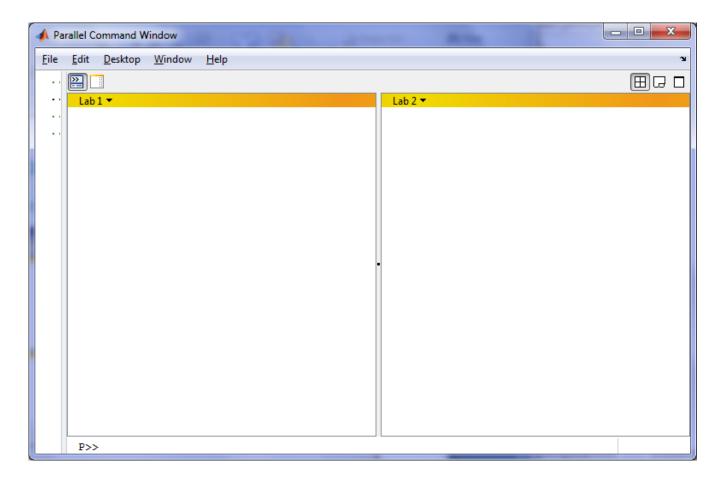
• profile viewer

Function listing

time	e calls	line
		1 function [time1,time2,time3] = newSPMDVersion
	1	2 n=10^8; step = 1/n;
		3 % Serial Version
	1	4 tic; s=0;
< 0.01	1	5 for i=1:n-1
8.24	99999999	<pre>6 x=(i-0.5)*step; s=s+4./(1+x^2);</pre>
10.49	999999999	7 end
	1	8 time1 = toc;
		9
		10 % Parallel Version
6.08	1	11 matlabpool open 4
	1	12 tic;
1.07	1	13 spmd
		14 slocal = myTestSum(n, step);
		15 end
	1	16 time2 = toc;
		17 % Parallel Version parfor
	1	18 nstep=10^8; step = 1/nstep;
	1	19 tic;s=0;
1.04	1	20 parfor i=1:nstep-1
		21 x=(i-0.5)*step;
		22 s=s+4./(1+x^2);
		23 end
3.32	1	24 time3 = toc;matlabpool close
	1	25 end

Parallel mode-II: pmode

>> pmode start



P>> pmode exit

pmode demo

P>> help magic % ask for help on a function

P>> PI = pi % set a variable on all the labs

P>> myid = labindex % lab ID

P>> all = numlabs % total No. of labs

P>> segment = [1 2; 3 4; 5 6] % create a replicated array on all the labs

P>> segment = segment + 10*labindex % perform on different labs

P >> x = magic(4) % replicated on every lab P >> y=codistirbuted(x) % partitioned among the lab

 $P >> z = y + 10^{*}$ labindex % operate on the distributed array whole

P>> combined = gather(y) % entire array in one workspace

The combined is now a 4-by-4 array in the client workspace. whos combined To see the array, type its name. combined

Demo: distributed array operations (repeat)

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- I	Lab 1					Lab 2	•			
•										
•										
						-				
	P>>	a=rand(4,4)				P>>	a=rand(4,4)			
		a =					a =			
		a –					a –			
		0.9173	0.4612	0.2155	0.4621		0.2951	0.7010	0.9143	0.7375
		0.6839	0.1562	0.4978	0.9846		0.0990	0.3821	0.2740	0.5407
			0.4626	0.2904	0.9587		0.3277	0.9602	0.6484	0.6348
		0.4809	0.8009	0.9071	0.5795		0.6902	0.7780	0.2781	0.0948
	P>>									

📣 Parallel (Command Window		-	-	-	-		<u> </u>
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^ 🖹	-							
Lab 1	•					Lab 2	2 -	
P>>	b=rand(2,8,cod	istributo	or())			P>>	<pre>b=rand(2,8,codistributor())</pre>	
I	This lab store	s b(:,1:4	ł).				This lab stores b(:,5:8).	
1	LocalP Codistribu	art: [2x4 tor: [1x1	-	butor1d]			LocalPart: [2x4 double] Codistributor: [1x1 codistributor1d]	
	getLocalPart(b)				P>>	> getLocalPart(b)	
	ans =						ans =	
	0.4267 0.5679			0.4703 0.5058	4		0.9787 0.1247 0.0250 0.1717 0.4276 0.9818 0.6535 0.4670	4 111
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▼ P>>								

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I	Lab 1	•						Lab 2	•					
I	P>>	gather(b)					^	P>>	gather(b)					
		ans =							ans =					
Í		Columns 1	through 7						Columns 1	through 7				
		0.4267	0.8379	0.6662	0.4703	0.9787			0.4267	0.8379	0.6662	0.4703	0.9787	
		0.5679	0.5489	0.7862	0.5058	0.4276			0.5679	0.5489	0.7862	0.5058	0.4276	
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	Lab 1	•					Lab 2	•				
		0.4267	0.8379	0.6662	0.4703			0.9787	0.1247	0.0250	0.1717	
		0.5679	0.5489	0.7862	0.5058			0.4276	0.9818	0.6535	0.4670	
	P>>	c=b'					P>>	c=b'				
		This lab sto	xoa a (1.4	• •				This lab sto	mag. g (E. 9	• •		
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		Codistri	butor: [1x	l codistri	.butor1d]	ŀ	•	Codistri	butor: [1x	1 codistri	ibutor1d]	
	P>>	getLocalPart	(C)				P>>	getLocalPart	(c)			
		ans =						ans =				
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		0.8379							0.9818			
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•						F I	•	III				Þ
-	P>>											

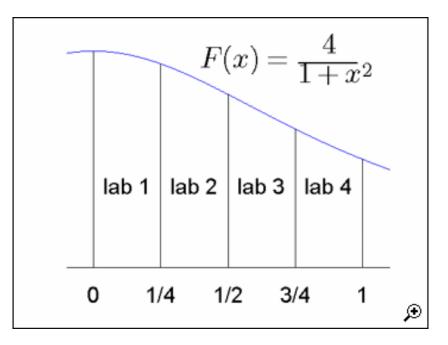
Parallel pi in pmode

use the fact that

$$\int_0^1 \frac{4}{1+x^2} dx = 4(atan(1) - atan(0)) = \pi$$

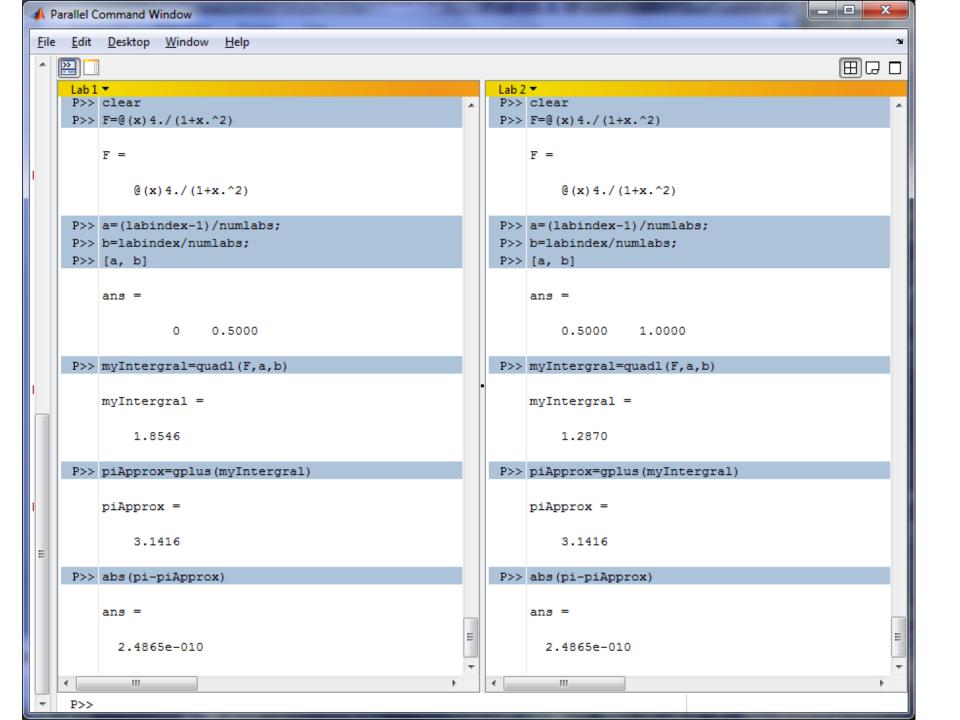
to approximate pi by approximating the integral on the left.

divide the work between the labs by having each lab calculate the integral the function over a subinterval of [0, 1] as shown in the picture



Steps

- All labs/workers will compute the same function: F=4/(1+x^2)
- Each worker/lab will calculate over a subinterval [a,b] of [0, 1], for 2 labs, the subinterval will be:
 - [0, 0.50] [0.50, 1.0] a = (labindex-1)/numlabs
 - b = labindex/numlabs
- Use a MATLAB quadrature method to compute the integral myIntegral = quadl(F, a, b)
- Add together to form the entire integral over [0,1] piApprox = gplus(myIntegral)

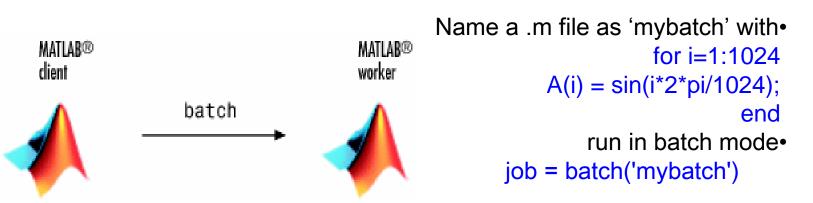


Parallel pi in matlabpool-mode

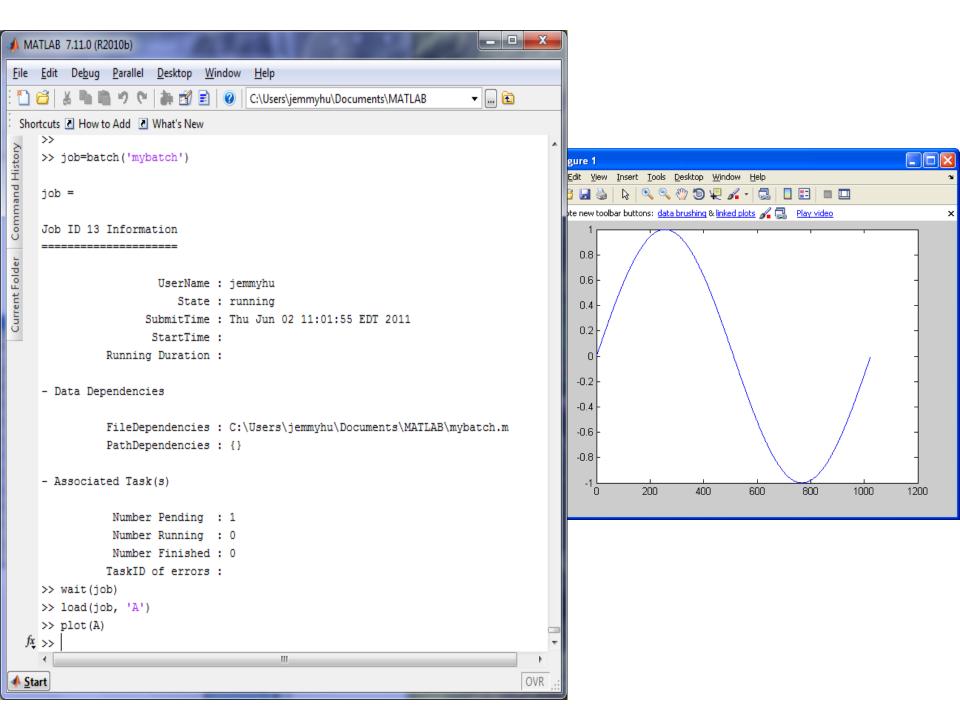
File Edit Text Go Cell Tgols Debug Desktop Window Help Image: Second							
<pre>************************************</pre>							
<pre>************************************</pre>							
<pre>(*) This file uses Cell Mode. For information, see the rapid code iteration video, the publishing video, or help.</pre>							
<pre>1</pre>							
<pre>2 % run Pi is spmd and matlabpool mode 3 % 4 % open matlabpool 5 = matlabpool open 6 7 % define the interval 8 = 0 spmd 9 = a = (labindex - 1)/numlabs; 10 = b = labindex/numlabs; 11 = fprintf('Subinterval: [%-4g, %-4g]\n', a, b); 12 = end 13 % labs use a quadrature method to approximate each integral 15 = 0 spmd 16 = myIntegral = quadl(@pctdemo_quadpi, a, b); 17 = fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 = end 19 % add the results together</pre>							
<pre>3</pre>							
<pre>4 %% open matlabpool matlabpool open 6 7 %% define the interval 8 spmd 9 -</pre>							
<pre>5 - matlabpool open 6 7 %% define the interval 8 spmd 9 -</pre>							
<pre>6 7 %* define the interval 8 spmd 9 -</pre>							
<pre>8 - spmd 9 -</pre>							
<pre>9 - a = (labindex - 1)/numlabs; b = labindex/numlabs; fprintf('Subinterval: [%-4g, %-4g]\n', a, b); 12 - end 13 14 %% labs use a quadrature method to approximate each integral 15 - spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
<pre>10 - b = labindex/numlabs; 11 - fprintf('Subinterval: [%-4g, %-4g]\n', a, b); 12 - end 13 14 %% labs use a quadrature method to approximate each integral 15 - spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
<pre>11 - fprintf('Subinterval: [%-4g, %-4g]\n', a, b); 12 - end 13 14 %% labs use a quadrature method to approximate each integral 15 - spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
<pre>12 - end 13 14 %% labs use a quadrature method to approximate each integral 15 spmd 16 myIntegral = quadl(@pctdemo_quadpi, a, b); 17 fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
<pre>13 14 13 14 %* labs use a quadrature method to approximate each integral 15 spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [*-4g, *-4g] Integral: *4g\n', a, b, myIntegral); 18 - end 19 20 ** add the results together</pre>							
<pre>14 %% labs use a quadrature method to approximate each integral 15 spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
<pre>15 spmd 16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 end 19 20 %% add the results together</pre>							
<pre>16 - myIntegral = quadl(@pctdemo_quadpi, a, b); 17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - end 19 20 %% add the results together</pre>							
17 - fprintf('Subinterval: [%-4g, %-4g] Integral: %4g\n', a, b, myIntegral); 18 - -end 19							
18 - - end 19 - 20 %% add the results together							
20 8% add the results together							
21 - Gand							
<pre>22 - piApprox = gplus(myIntegral);</pre>							
23 - fprintf('piApprox: %-4g\n', piApprox);							
24 end 25							
%% results in client							
<pre>** results in client approx1 = piApprox{1}; % 1st element holds value on lab 1.</pre>							
<pre>27 - approx1 = piApprox{1}; % 1st element holds value on lab 1. 28 - fprintf('pi : %.18f\n', pi);</pre>							
<pre>29 - fprintf('Approximation: %.18f\n', approx1);</pre>							
30 - fprintf('Error : %g\n', abs(pi - approx1));							
31							
32 %% close matlabpool							
33 - matlabpool close -							
pctdemo Ln 5 Col 16 OVR							

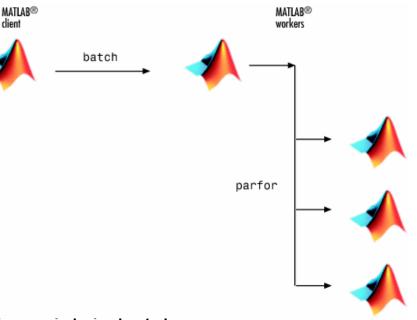
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Shortcuts 🔄 How to Add 💽 What's New
     >> pmode start
Current Folder Command History
    Starting pmode using the 'local' configuration ... connected to 2 labs.
    >> pmode exit
     Sending a stop signal to all the labs ... stopped.
     >> run('C:\Users\jemmyhu\Documents\MATLAB\pctdemo.m')
     Starting matlabpool using the 'local' configuration ... connected to 2 labs.
    Lab 1:
       Subinterval: [0 , 0.5 ]
    Lab 2:
      Subinterval: [0.5 , 1 ]
    Lab 1:
       Subinterval: [0 , 0.5 ] Integral: 1.85459
    Lab 2:
       Subinterval: [0.5, 1] Integral: 1.287
    Lab 1:
      piApprox: 3.14159
     Lab 2:
      piApprox: 3.14159
    pi
        : 3.141592653589793100
    Approximation: 3.141592653838447500
                  : 2.48654e-010
     Error
     Sending a stop signal to all the labs ... stopped.
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Batch mode



- The batch command does not block MATLAB, so you must wait for the job to finish before you can retrieve and view its results: wait(job)
- The load command transfers variables from the workspace of the worker to the workspace of the client, where you can view the results:
 - load(job, 'A') plot(A)
 - When the job is complete, permanently remove its data:• destroy(job)





A batch parallel loop

% mybatch parfor i=1:1024 A(i) = sin(i*2*pi/1024); end

job = batch('mybatch', 'configuration, 'local', 'matlabpool', 2)

% To view the results: wait(job) load(job, 'A') plot(A)

% remove its data: destroy(job)

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Shortcuts 🖪 How to Add 💽 What's New	\vdots	
	1 %run in batch mode	
job =	2 - job = batch('pmybatch')	
Job ID 22 Information	3 %The batch command does not block MATLAB, so you must wait 4 %for the job to finish before you can retrieve and view its results:	
	<pre>5 - wait(job)</pre>	
	6 %The load command transfers variables from the workspace of the	
UserName : elias	7 %worker to the workspace of the client, where you can view the	
State : running	8 %results:	
SubmitTime : Fri Nov 23 07:23:52 EET 2012	9 - load(job, 'A')	
StartTime :	<pre>10 - plot(A) 11 %When the job is complete, permanently remove its data:</pre>	
Running Duration :	12 - destroy(job)	
- Data Dependencies	13	
	Figure 1	
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PathDependencies : {}		
- Associated Task(s)		
Number Pending : 1		
Number Running : 0	0.8-	
Number Finished : 0		
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Current Folder Command History
     >> job=batch('mybatch par', 'configuration', 'local', 'matlabpool', 1)
     job =
     MatlabPool Job ID 17 Information
         _____
                       UserName : jemmyhu
                           State : queued
                     SubmitTime : Thu Jun 02 11:09:39 EDT 2011
                      StartTime :
               Running Duration :
     - Data Dependencies
               FileDependencies : C:\Users\jemmyhu\Documents\MATLAB\mybatch par.m
               PathDependencies : {}

    Associated Task(s)

                Number Pending : 2
                Number Running : 0
                Number Finished : 0
               TaskID of errors :
     - Scheduler Dependent (MatlabPool Job)
         MaximumNumberOfWorkers : 2
         MinimumNumberOfWorkers : 2
     >> wait(job)
     >> load(job, 'A')
     >> plot(A)
     >> destroy(job)
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Start
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                                                                                                     ×
      function interactiveToBatch()
 1
 2
         % interactive to batch
 3
 4
         %% for loop
 5 -
         A = cell(1,4);
 6 -
      for(i=1:4)
 7 -
         A{i} = magic(i);
 8 -
         end
         A{:}
 9 -
10
11
         %% interactive parallel for loop
         A = cell(1,4);
12 -
      parfor(i=1:4)
13 -
14 -
         A(i) = magic(i);
15 -
        end
         A{:}
16 -
17
18
         %% batch job
         jm = findResource('scheduler', 'configuration', 'local');
19 -
20 -
         job = createJob(jm);
       for(i=1:4)
21 -
22 -
             createTask(job, @magic, 1, {i});
23 -
        end
24 -
        submit(job);
        waitForState(job);
25 -
        A = getAllOutputArguments(job);
26 -
27 -
         A{:}
28 -
         end
                                                     interactiveToBatch
                                                                               Ln 1
                                                                                        Col 1
```

Key Function List

Job Creation

<u>createJob</u> Create job object in scheduler and client <u>createTask</u> Create new task in job <u>dfeval</u> Evaluate function using cluster

Interlab Communication Within a Parallel Job
 labBarrier Block execution until all labs reach this call
 labBroadcast Send data to all labs or receive data sent to all labs
 labindex Index of this lab
 labReceive Receive data from another lab
 labSend Send data to another lab
 numlabs Total number of labs operating in parallel on current job

Job Management

<u>cancel</u> Cancel job or task <u>destroy</u> Remove job or task object from parent and memory <u>getAllOutputArguments</u> Output arguments from evaluation of all tasks in job object <u>submit</u> Queue job in scheduler wait Wait for job to finish or change states

Typical Use Cases

• Parallel for-Loops (parfor)

allowing several MATLAB workers to execute individual loop iterations simultaneously

restriction on parallel loops is that no iterations be allowed to depend on any other iterations.

Large Data Sets

allows you to distribute that array among multiple MATLAB workers, so that each worker contains only a part of the array

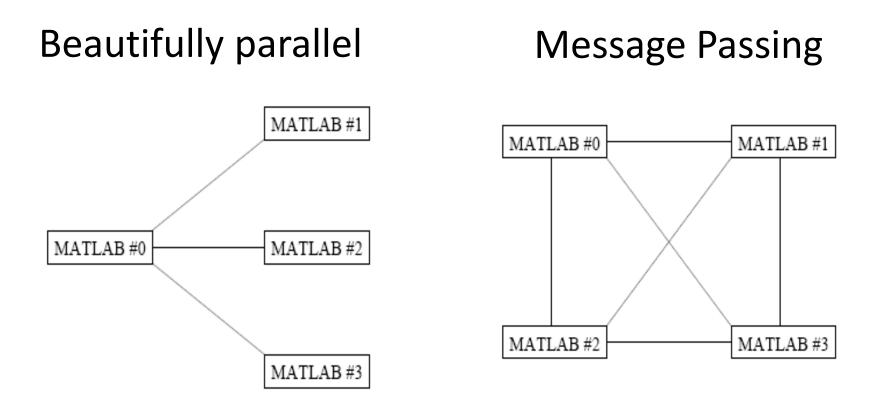
Each worker operates only on its part of the array, and workers automatically transfer data between themselves when necessary

• Batch Jobs

offload work to a MATLAB worker session to run as a batch job.

the MATLAB worker can run either on the same machine as the client, or if using MATLAB Distributed Computing Server, on a remote cluster machine.

Matlab in Parallel



e.g., Multi, paralize, Plab, ParMatlab

e.g., MultiMatlab, CMTM, DPToolbox,MatlabMPI, pMatlab

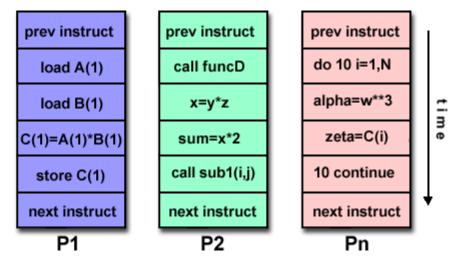
www-math.mit.edu/~edelman/homepage/papers/pmatlab.pdf

MPI+SMP Parallel Paradigms

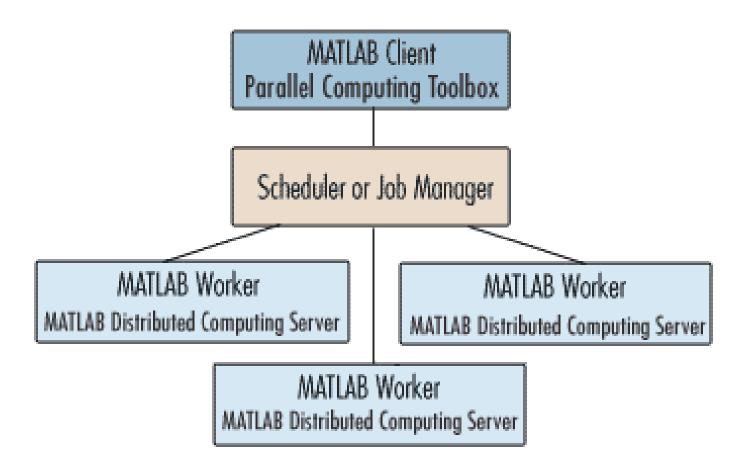
<u>http://ist.uwaterloo.ca/ew/saw/parallel/FLAS</u>
 <u>H/swf/mpi+smp.swf</u>

MIMD

- Currently, most common type of parallel computer
- Every processor may be executing a different instruction stream
- Every processor may be working with a different data stream
- Execution can be synchronous or asynchronous, deterministic or nondeterministic
- Examples: most current supercomputers, networked parallel computer clusters and "grids", multi-processor SMP computers, multi-core PCs.
- Note: many MIMD architectures also include SIMD execution sub-components



Parallel Computing Toolbox and MATLAB Distributed Computing

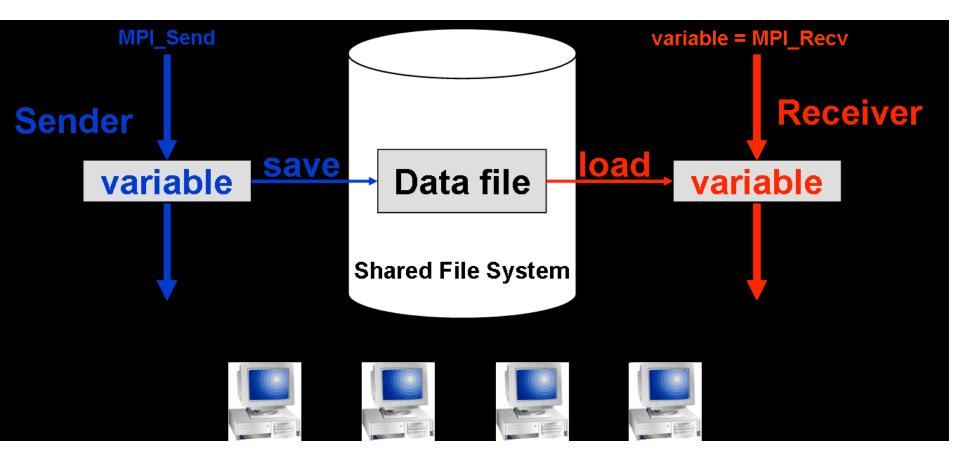


MPI Library

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Shortcuts 🛛 How to Add 💽 What's New	
Compute Time (sec) Launch+Comm Time (sec) >> help MPI Contents of mpi:	= 2.7435 = 0.0068936
dctProfStripAnchors	- Removes anchors that evaluate MATLAB code from
labBarrier	- block until all labs have entered the barrier
labBroadcast	- send data to all labs
labProbe	 test to see if messages are ready to labReceive
labReceive	- receive data from another lab
labSend	- send data to another lab
labSendReceive	- simultaneously send to and receive from other labs
labindex	- return the ID for this lab
mpiFinalize	- perform MPI finalization
mpiInit	- perform MPI initialization
<u>mpiInitialized</u>	- query whether MPI is initialized
mpiLibConf	- return the location of an MPI implementation
mpiParallelSessionEnding	- called at the end of a parallel session
mpiParallelSessionStarting	- called at the start of a parallel session
mpiSettings	- set various options for MPI communication
mpigateway	- gateway to internal MPI functionality
mpiprofile	- Profile parallel communication and execution times.
mpiprofview	- Displays the html profiler interface or returns the html
mpiprofviewgateway	- PROFVIEWGATEWAY Profiler HTML gateway function.
numlabs	- return the total number of labs operating in parallel

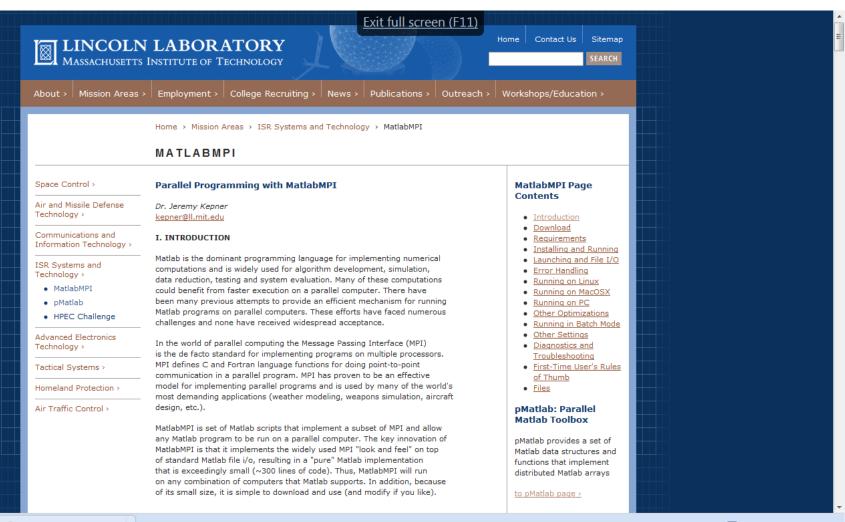
fx >> |

MatlabMPI implements the fundamental communication operations in MPI using MATLAB's file I/O functions.

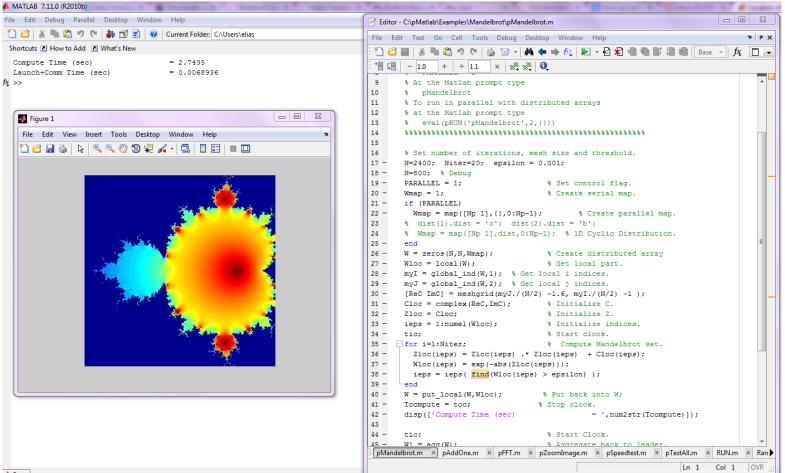


MatlabMPI

http://www.ll.mit.edu/mission/isr/matlabmpi/matlabmpi.html#introduction



pMatLab example



FFT with pMatLab

MATLAB 7.11.0 (R2010b) C:\pMatlab\Examples\FFT\pFFT.m				Sunction details for pFFT									
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<pre>>> profile on Np Pid Distributed vector size (words) Distributed vector size (bytes) Local vector size (bytes)</pre>	= 1 = 0 = 1048576 = 16777216 = 16777216	1 2 3 4 5 6	<pre>% This script implements % the most common way of % contains a parallel imp % % Parameters:</pre>	a simple 1D FFT benchmark us implementing a 1D FFT in par plementation of this algorith	Concentrated 25-Nov-2012 21:05:34 using cpu time. Script in file C:\pMatlab\Examples\FFT\pFT.m Copy to new window for comparing multiple runs Parents (calling functions)								
Allocation Time (sec) Launch Time (sec)	= 0.11417 = 0.01701	8	7 % * PARALLEL - Enable the pMatlab library.			No parent							
Launch Time (sec) = 0.01701 Begin 1st CornerTurn			9 % * VALIDATE - Enable validation of FFT result.			Lines where the most time was spent							
Begin FFT of 2nd Dimension		10	8		Line Number	Code			Calls	Total Time	% Time	Time Plot	
Begin 2nd CornerTurn Compute time (sec)	= 0.13974	11 12	<pre>% * ERROR_LIMIT - Error %</pre>	r limit used in validation.	86	Xloc = fft(Xlo	oc,[],2);		1	0.123 s	27.3%		
Communication time (sec)	= 0.15662	13	-	or to FFT, must be divisible 1	45	X = complex(ra	and(1.N.Xmap)	.ran	1	0 077 s	17.1%		
Run time (sec)	= 0.29636	14 15	<pre>%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%</pre>	81	<pre>X = transpose grid(X);</pre>				0.063 s	14.0%	-		
Performance (Gigaflops) Bandwidth (Gigabytes/sec)	= 0.35382 = 0.32137	16	<pre>% PARALLEL = 0</pre>	it distributed arrays, set									
>> profile viewer		17	% At the Matlab prompt ty	ype	92	X = transpose_	grid(X);		1	0.043 s	9.5%		
$f_{x} >>$		18 19	<pre>% pFFT % To run in serial with c</pre>	distributed arrays set	100	X = transpose_	grid(X);		1	0.041 s	9.1%	•	
		20	<pre>% PARALLEL = 1</pre>	instituteu allays, set	All other lines					0.104 s	23.1%		
		21	<pre>% At the Matlab prompt ty</pre>	/pe	Totals					0.451 s	100%		
		22 23 24	<pre>% pFFT % To run in parallel with (% at the Matlab prompt type % eval(pRUN('pFFT',2,{))) %**********************************</pre>	<i>T</i> pe	Children (called functions)								
		25 26 27 28 - 29 -			Function Name		Function Type	Calls T	otal Time	% Time	Time Plot		
					transpose_grid		function	3 0	.147 s	32.6%			
							function	2 0	.065 s	14.4%			
		30 -	$N = 2^{20*Np}; M = N/Np;$ N = 2^20*Np; M = N/Np;	<pre>% Debug.</pre>			function	12 0	.012 s	2.7%	1		
		31 - 32 -	VALIDATE = 0; ErrorRate = eps;	% Turn validation on or of % Set error to machine pre-	dmat.agg		function	1 0	.011 s	2.4%	1		
		32 - Errorkate =	Errorkate - eps;				function		010 s	2.2%	1	-	
		34 -	Xmap = 1;	<pre>% Serial map.</pre>	· · · · ·		function		.010 s	1.3%		-	
		35 - 36 -	<pre>if PARALLEL Xmap = map([1 Np],{},0:</pre>	Np-1); % Parallel map.	map.zeros							-	
		•			map.map		function		.003 s	0.7%		-	
▲ Start		: pMande	pMandelbrot.m × pAddOne.m × pFFT.m × pZoomImage.m × pSpeedtest.m			Pid fu		2 0	.001 s	0.2%			
				script	•								

Available examples:

- xbasic.m Extremely simple MatlabMPI program that prints out the rank of each processor.
- basic.m Simple MatlabMPI program that sends data from processor 1 to processor 0.
- multi_basic.m Simple MatlabMPI program that sends data from processor 1 to processor 0 a few times.
- probe.m Simple MatlabMPI program that demonstrates the using MPI_Probe to check for incoming messages.

broadcast.m Tests MatlabMPI broadcast command. basic_app.m Examples of the most common usages of MatlabMPI. basic_app2.m Examples of the most common usages of MatlabMPI. basic_app3.m Examples of the most common usages of MatlabMPI. basic_app4.m Examples of the most common usages of MatlabMPI. MatlabMPI test parallel image processing application. blurimage.m speedtest.m Times MatlabMPI for a variety of messages. synch_start.m Function for synchronizing starts. machines.m Example script for creating a machine description. Wrapper for using an example as a unit test. unit_test.m unit_test_all.m Calls all of the examples as way of testing the entire library. unit_test_mcc.m Wrapper for using an example as a mcc unit test.

unit_test_all_mcc.m Calls all of the examples using MPI_cc

as way of testing the entire library.

xbasic

```
% Basic Matlab MPI script that
% prints out a rank.
%
% To run, start Matlab and type:
%
%
  eval( MPI_Run('xbasic',2,{}) );
%
 Or, to run a different machine type:
%
%
%
  eval( MPI_Run('xbasic',2,{'machine1' 'machine2'}) );
%
% Output will be piped into two files:
%
%
  MatMPI/xbasic.0.out
%
  MatMPI/xbasic.1.out
%
% MatlabMPI
% Dr. Jeremy Kepner
% MIT Lincoln Laboratory
% kepner@ll.mit.edu
```

```
% Initialize MPI.
MPI Init;
% Create communicator.
comm = MPI_COMM_WORLD;
% Modify common directory from default for better performance.
% comm = MatMPI Comm dir(comm,'/tmp');
% Get size and rank.
comm_size = MPI_Comm_size(comm);
my_rank = MPI_Comm_rank(comm);
% Print rank.
disp(['my rank: ',num2str(my rank)]);
% Wait momentarily.
pause(2.0);
% Finalize Matlab MPI.
MPI Finalize;
disp('SUCCESS');
if (my_rank ~= MatMPI_Host_rank(comm))
  exit;
end
```

Demo folder ~/matlab/, watch top at the other machine

```
🖉 vdwarf2.ee.bgu.ac.il - PuTTY
vdwarf2.ee.bgu.ac.il> matlab -nodesktop -nodisplay -nojvm
                               < M A T L A B >
                   Copyright 1984-2007 The MathWorks, Inc.
                          Version 7.5.0.338 (R2007b)
                                August 9, 2007
        Your MATLAB license will expire in 11 days.
        Please contact your system administrator or
        The MathWorks to renew this license.
  To get started, type one of these: helpwin, helpdesk, or demo.
  For product information, visit www.mathworks.com.
>> eval( MPI Run('xbasic', 2,{'vdwarf3','vdwarf4'}) );
Launching MPI rank: 1 on: vdwarf4
Launching MPI rank: 0 on: vdwarf3
unix launch =
 rsh vdwarf4 -n 'cd /users/agnon/misc/tel-zur/matlab; /bin/sh ./MatMPI/Unix Comm
ands.vdwarf4.1.sh &' &
 rsh vdwarf3 -n 'cd /users/agnon/misc/tel-zur/matlab; /bin/sh ./MatMPI/Unix Comm
ands.vdwarf3.0.sh &' &
* * * * * * * *
```

Parallel Matlab (Octave) using pMatlab

Global arrays – "...Communication is hidden from the programmer; arrays are automatically redistributed when necessary, without the knowledge of the programmer..."

"...The ultimate goal of pMatlab is to move beyond basic messaging (and its inherent programming complexity) towards higher level parallel data structures and functions, allowing any MATLAB user to parallelize their existing program by simply changing and adding a few lines,

Source: http://www.ll.mit.edu/mission/isr/pmatlab/pMatlab_intro.pdf

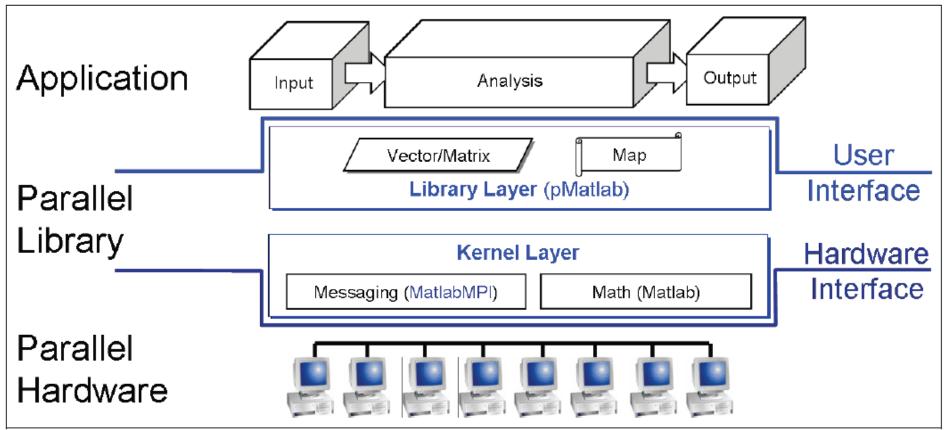
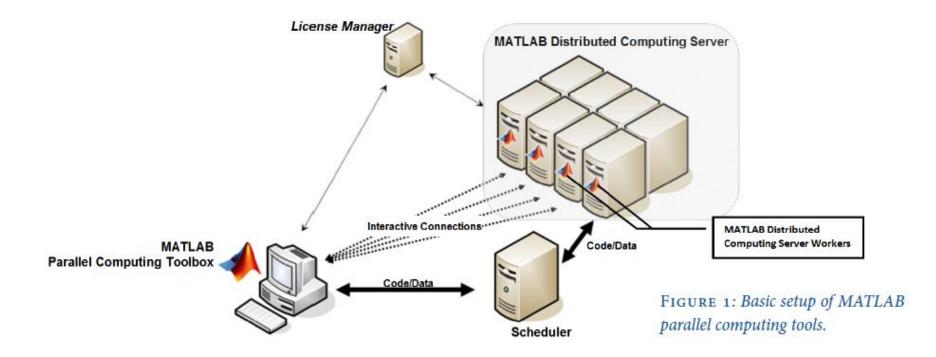
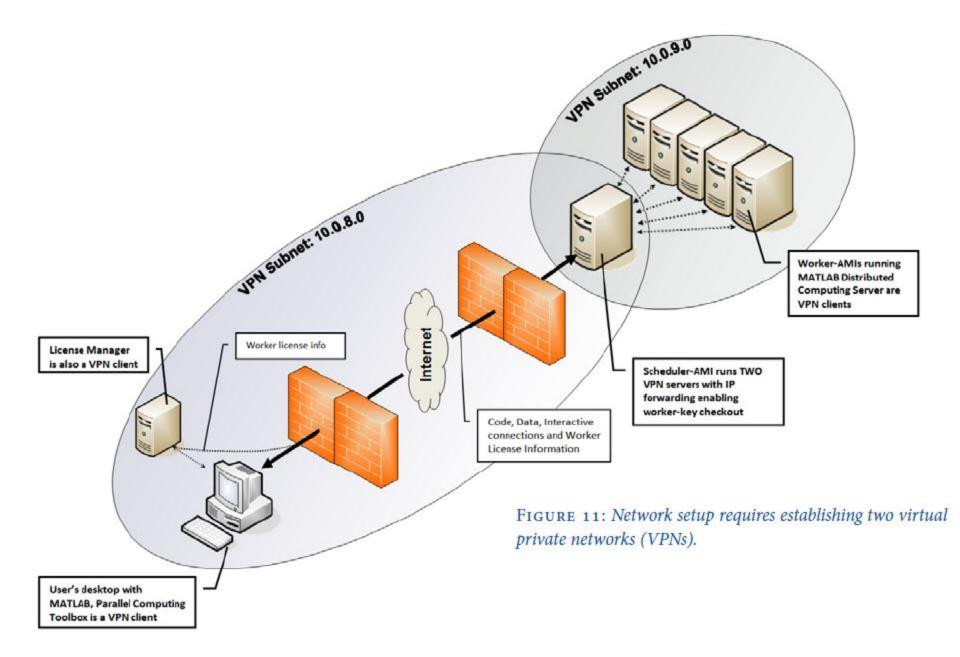


Figure 11 – Parallel MATLAB consists of two layers. pMatlab provides parallel data structures and library functions. MatlabMPI provides messaging capability.

Parallel Computing with Matlab on Amazon Cloud

MATLAB Parallel Computing Tools: Basic Setup and Requirements





Example - Problem Description

- System of 2^6 (=64) square matrices
 - − Each matrix → Sparse, square,2^17 (=131072) dimension
- To extract first 100 eigenvectors
 'eigs' function is used
- See handout for the code
- Each matrix calculation is distributed

Example - Serial Matlab

- 'eigen' is a function
 - Input : (vector of random numbers, dimension of the matrix)
 - Output : eigenvectors

$$n = 2^{16};$$

$$p = 2^{7};$$

$$e = rand (n, 1, p);$$

$$for i = 1 : p$$

$$a = e(:, :, i);$$

$$ans(i) = eigen(a, n);$$

$$end$$

Example-Parallel Matlab

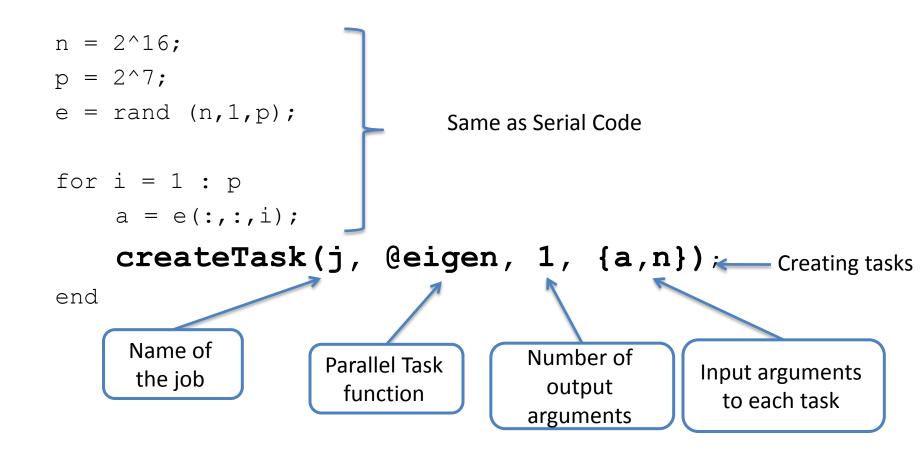
1. Find available distributed computing resources
 (findResource function)

```
nprocs = [ getenv('DMATLAB_NPROCS') ]
np = sscanf( nprocs, '%d' )
mgr_name = [ getenv('JOBMANAGER') ]
mgr_host = [ getenv('JOBMANAGERHOST') ]
jm =findResource('jobmanager', 'Name', mgr_name, 'LookupURL', mgr_host);
```

2. Create distributed job

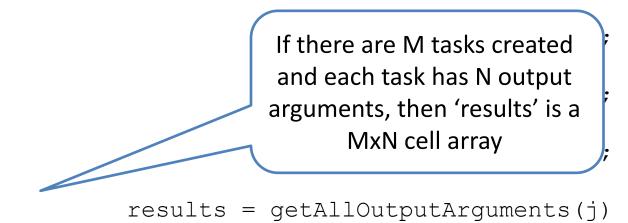
Example – Parallel Matlab

3. Create Tasks for each worker



Example – Parallel Matlab

4. Submit the job and wait for the results



5. Remove the individual task or parent job object

destroy(j);