Some changes in snow and R

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- Some new ideas for me:
 - Grid computing
 - MapReduce
 - OpenMP
- Some opportunities:
 - Rethink some aspects of snow design
 - Make progress on parallel vectorized arithmetic for R



- better error handling
- integrating load balancing into all functions
- R-level collection of timing information
- non-parallel testing framework
- persistent data on nodes
- limited inter-node communication
- sensible handling of user interrupts



- Errors in clusterXYZ functions used to be returned as try-error objects.
- New version will signal an error on the master if there is an error on any node.
- This is better but not always ideal: Sometimes one good result is enough.



- Originally clusterApply required at most as many elements as cluster nodes.
- Longer lists can be handled by clusterApplyLB
- This does not provide a deterministic option for longer vectors.
- It also leaves out the new clusterMap function.
- The new version allows longer vectors.
- By default nodes are recycled.
- This produces deterministic job/node assignments.



- Load balancing might be useful with functions other than just clusterApply.
- Motivated by OpenMP, one option is to
 - allow a SCHEDULE argument with values "static" or "dynamic" got all functions
 - a variant is to allow a boolean LoadBalance argument
 - have the parXYZ functions take a ChunkSize argument
- Default will be static.
- Open question: can parallel RNG streams be tied to jobs in a simple way so results from with load-balancing can be reproducible?

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- xpvm is very useful, but requires pvm.
- xmpi is similar but restricted to LAM.
- An alternative is to collect timing information in R:
 - In the master record the start and finish of each send/recv.
 - In the nodes, record duration of computation and send back with result.
- Need to decide interface for collecting the data.
- One possibility, motivated by **Rprof**:
 - traceCluster(file=''foo.trace'') to start recording to a file
 - traceCluster(NULL) to stop recording
- Then need some functions to read trace file and produce graphs.
- Will experiment with this in the next month or so.



• It may be useful to have a "null cluster" so that

cl <- makeNULLcluster(4)
clusterApply(cl, ...)</pre>

works within the master process

- This will help with
 - debugging
 - running small jobs without parallel complications
- Some detail issues:
 - should the cluster size argument matter?
 - should random number streams behave as in the parallel version?



- It can be useful to leave large computed values on nodes for further computation.
- Global variables can be used but are awkward and not very clean.
- A better option may be to have a means of returning only a remote object reference.
- These remote objects can then be passed to subsequent calls.
- Once the master no longer has a reference to a remote object it can be garbage collected.

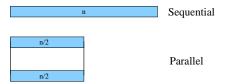


• A next step is to allow remote data to move between nodes, e.g.

- think of the nodes as arranged in a circle
- each node passes its data to the node to its left
- This leads to a model called Bulk Synchronous Parallel (BSP) computing.
- BSP has some interesting theoretical properties
 - a cost model for comparing parallel algorithms in terms of simple machine parameters
 - deadlock-free
- BSP has been used as the basis for parallel computing support for several high level languages.
- An initial (and maybe inefficient) BSP extension may be available soon.

• Basic idea for computing f(x[1:n]) on a two-processor system:

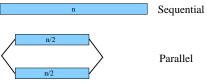
- Run two worker threads.
- Place half the computation on each thread.
- Ideally this would produce a two-fold speed up.





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• Reality is a bit different:

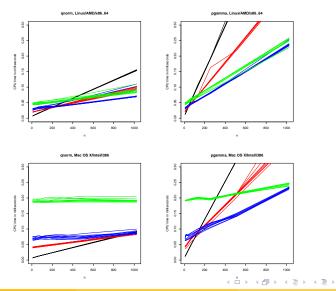


- There is synchronization overhead.
- Use of shared resources is sequential (memory, bus, ...)
- Parallelizing will only pay off if *n* is large enough.
 - For some functions, e.g. <code>qbeta</code>, $n \approx 10$ may be large enough.
 - For some, e.g. qnorm, $n \approx 1000$ is needed.
 - For basic arithmetic operations $n \approx 30000$ may be needed.
- Careful tuning to insure improvement will be needed.
- Some aspects will depend on architecture and OS.

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Parallelizing Vector Operations Some Experimental Results





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Some observations:

- Times are roughly linear in vector length.
- Intercepts on a given platform are roughly the same for all functions.
- If the slope for P processors is s_P , then at least for P = 2 and P = 4,

$$s_P \approx s_1/P$$

• Relative slopes of functions seem roughly independent of OS/architecture.

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• A simple strategy:

- Compute relative slopes once, or average across several setups.
- Base line is a single element dnorm computation.
- For each OS/architecture combination compute the intercepts.
- Estimate the values $N_2(f)$ such that using P = 2 is faster if $n > N_2(f)$.
- Use $N_4(f) = 2N_2(f)$ and $N_8(f) = 4N_2(f)$.

Some intercepts, in units of a single element <u>dnorm</u> computation:

- about 200 for Linux/AMD/x86_64
- about 500 for Mac OS X 10.4/Intel/i386
- between 300 and 400 for Win32/Intel(?)

Parallelizing Vector Operations



Some N2(f) Values on Linux

sqrt	o
sin	0
cos	0
exp	0
dnorm	0
pnorm	0
qnorm	0
dgamma	0
pgamma	0
qgamma	o
pbeta	••• • •
qbeta	• •
ptukey	• •
qtukey	0
	0 500 1000 1500 2000

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- Need to use threads
- One possibility: using raw pthreads
- Better choice: use Open MP
- Open MP consists of
 - compiler directives (#pragma statements in C)
 - a runtime support library
- Most commercial compilers support Open MP.
- gcc 4.2 supports Open MP.
- Redhat has back-ported Open MP into gcc 2.4.1 on RH, Fedora.
- MinGW also supports Open MP; an additional pthreads download is needed.

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• Basic loop for a one-argument function:

```
#pragma omp parallel for if (P > 1) num_threads(P) \
    default(shared) private(i) reduction(&&:naflag)
    for (i = 0; i < n; i++) {
        double ai = a[i];
        MATH1_LOOP_BODY(y[i], f(ai), ai, naflag);
    }
</pre>
```

• Steps in converting to Open MP:

- check f is thread-safe; modify if not
- rewrite loop to work with the Open MP directive
- test without Open MP, then enable Open MP

- Some things that are not thread-safe:
 - use of global variables
 - R memory allocation
 - signaling warnings and errors
 - user interrupt checking
 - creating internationalized messages (calls to gettext)
- Random number generation is also problematic.
- Functions in nmath that have not been parallelized yet:
 - Bessel functions
 - Wilcoxon, signed rank functions
 - random number generators



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• Package pnmath is available at

http://www.stat.uiowa.edu/~luke/R/experimental/

- This requires a version of gcc that
 - supports Open MP
 - allows dlopen to be used on libgomp.so

Our current systems don't satisfy this.

- A version using just pthreads is available in pnmath0. This should work on current R on our systems.
- Loading these packages replaces builtin operations by parallelized ones.
- For Linux, Mac OS X predetermined intercept calibrations are used.
- For other platforms a calibration test is run at package load time.
- The calibration can be run manually by calling calibratePnmath
- Hopefully we will be able to include this in R 2.7 or 2.8.

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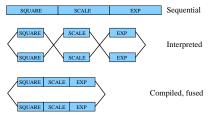
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Connection to Compilation

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- Developing a byte code compiler for R is an ongoing project.
- The codetools package is a by-product.
- Compilation will also be useful for parallelizing vector operations:
 - Many vector operations occur in compound expressions, like exp(-0.5*x²)
 - A compiler may be able to fuse these operations:



• Compilation may also allow many simple uses of apply functions and sweep to be parallelized.

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Challenges

• Tuning issues:

- Parallelization interface for package use.
- Extensible byte code for package use.
- Error handling and user interrupts.
- Performance may vary with inputs. Load balancing may be useful.
- Hardware/OS plays a role.
- Competing system usage may be important.

- Generic functions and non-default methods.
- Declarations may be useful.