

# Global HPCC Benchmarks in Chapel

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# Chapel in a nutshell

## *Chapel:*

- a new parallel language being developed by Cray Inc.
- part of DARPA's HPCS\* program
- first public release occurred this past weekend

\*HPCS = High Productivity Computing Systems

# When we last saw you at HPCC...

## HPCC 2006: Chapel “elegance only” entry

- **goal:** show where Chapel was headed
- 3 benchmarks: STREAM Triad, Random Access, FFT
- written with elegance and scalability in mind
- compiled and executed correctly, *but:*
  - only supported single-threaded execution
  - leaked memory
  - =>no performance

## This year's entry

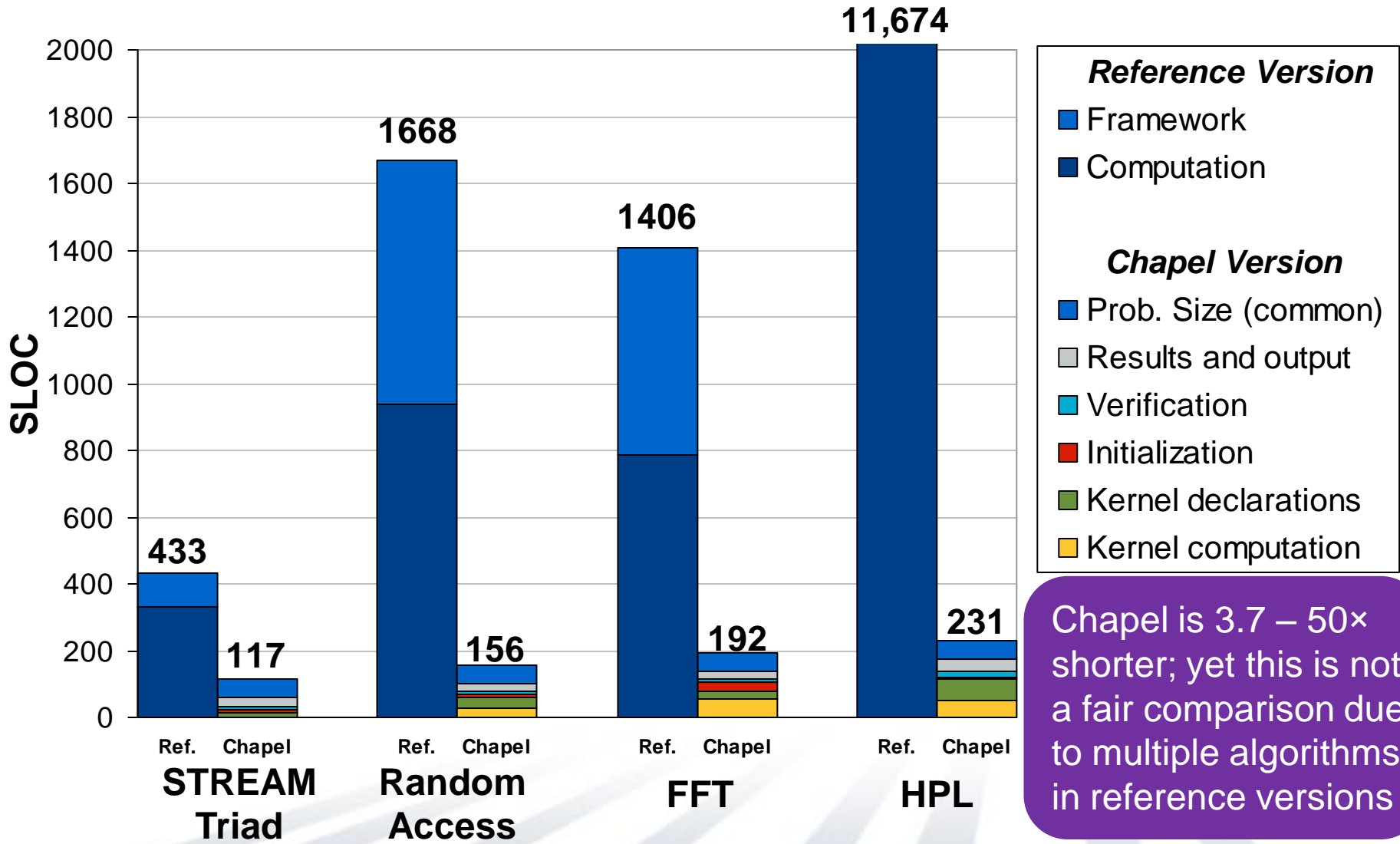
- First public performance numbers for Chapel execution
- First distributed memory execution of our data parallel features
- As intended, our code is quite similar to 2006 entry
- First locality-sensitive implementation of HPL in Chapel

## Please set your expectations appropriately:

- This is a snapshot of a work in progress, not the final word
- Our first *distribution* ran for the first time only two months ago

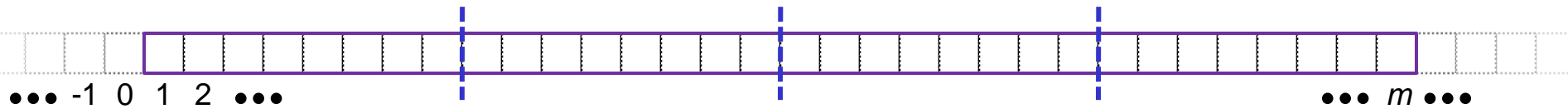
Focus less on our current performance  
and more on *how* we got it

# Code Size Summary (SLOC)

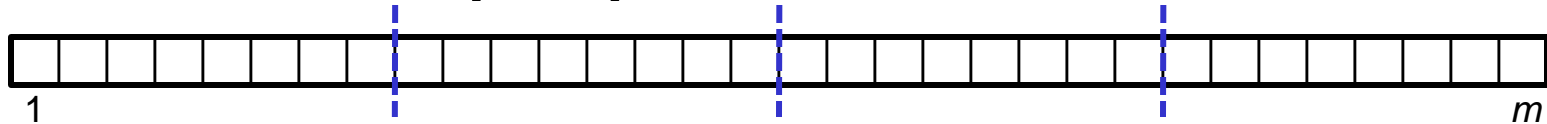


# STREAM Triad in Chapel

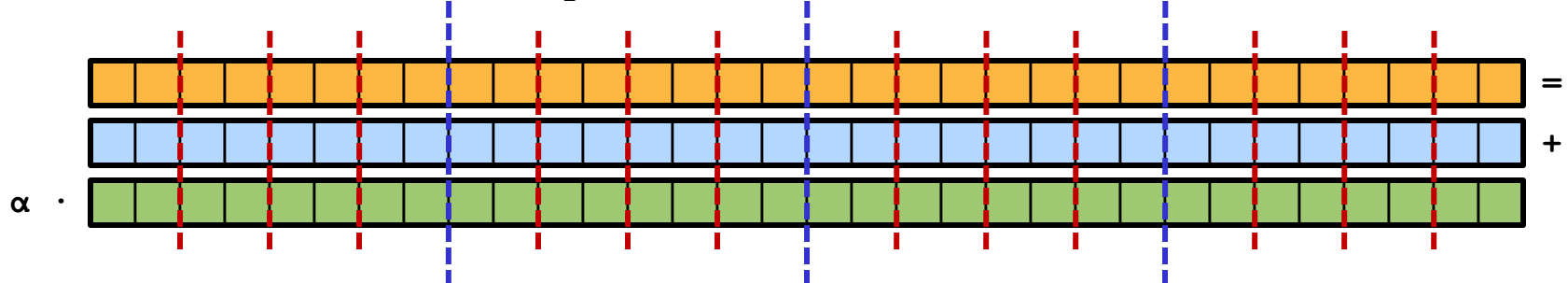
```
const BlockDist = new Block1D(bbox=[1..m], tasksPerLocale=...);
```



```
const ProblemSpace: domain(1, int(64)) distributedBlockDist
    = [1..m];
```



```
var A, B, C: [ProblemSpace] real;
```

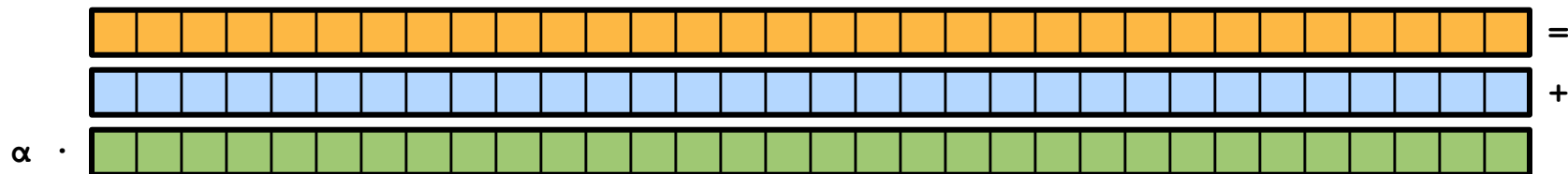


```
forall (a, b, c) in (A, B, C) do
    a = b + alpha * c;
```

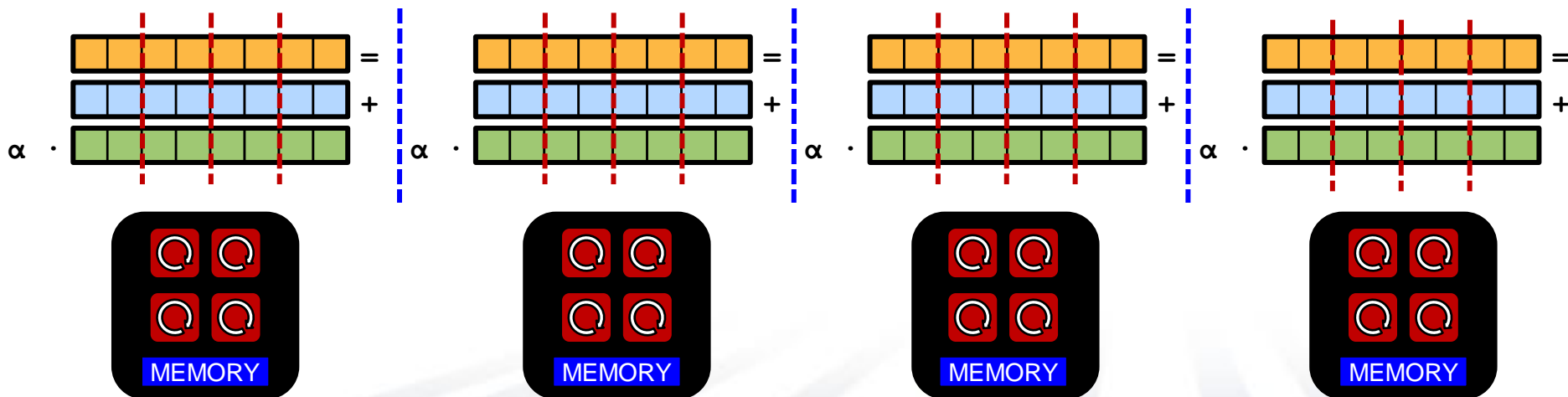
# Chapel Distributions

**Distributions:** “Recipes for parallel, distributed arrays”

- help the compiler map from the computation’s global view...



...down to the *fragmented*, per-processor implementation



# Chapel Distributions

- (Advanced) Programmers can write distributions in Chapel
- Chapel will support a standard library of distributions
  - *research goal*: using the same mechanism that users would
- Block1D is our first such distribution
  - *our compiler has no semantic knowledge of block distributions*
  - only of a distribution's interface--how to...
    - ...create domains and arrays using that distribution
    - ...map indices to locales
    - ...access array elements
    - ...iterate over indices/array elements
      - sequentially
      - in parallel
      - in parallel and zippered with other parallel iterable types
    - ...and so forth...

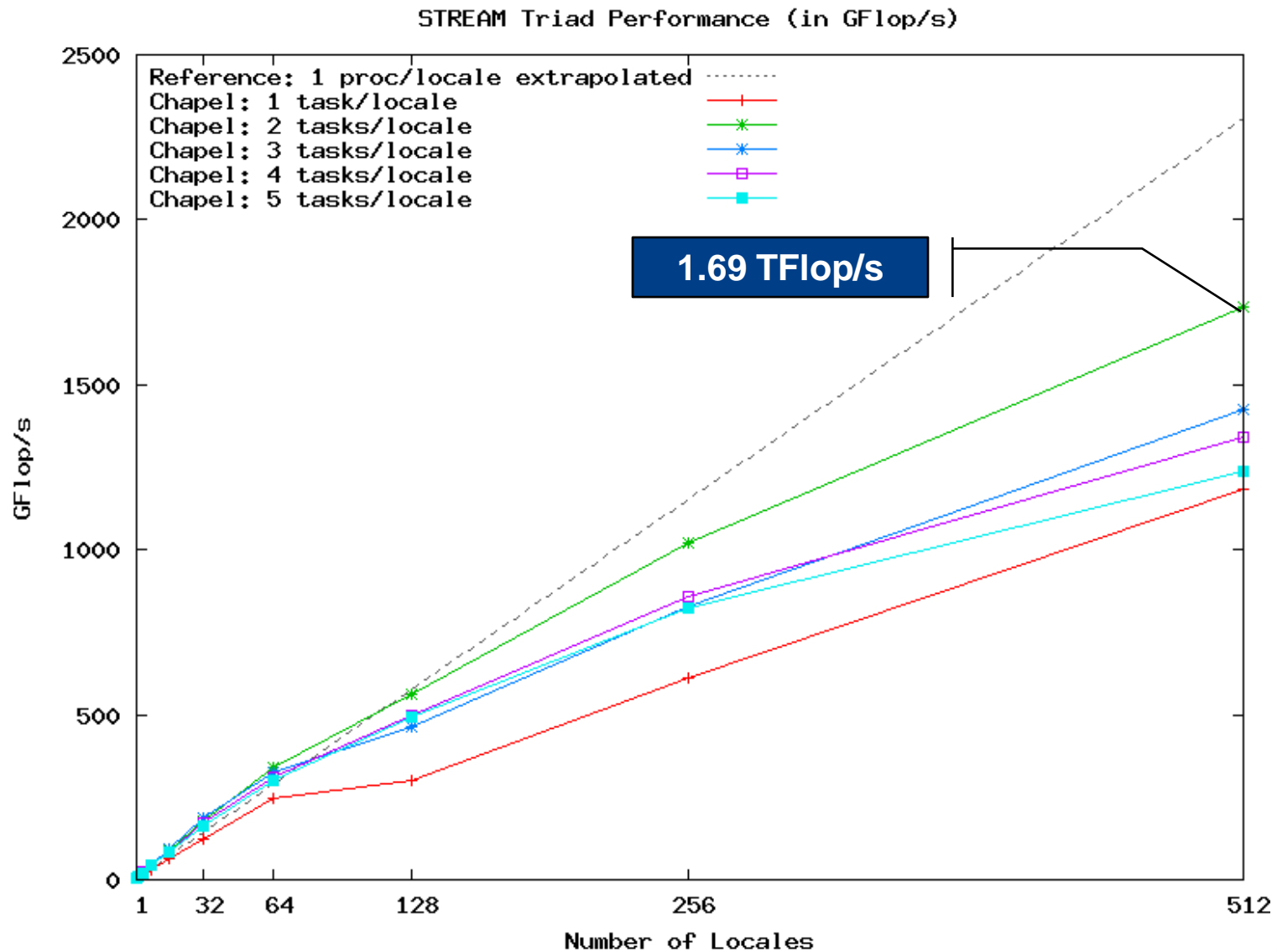


# Experimental Platform

<i>machine characteristic</i>	<i>value</i>
name	jaguar
model	Cray XT4
location	ORNL
# compute nodes	7,832
compute node processor	2.1 GHz AMD Opteron
cores per node	4
total user RAM per node	7.68 GB

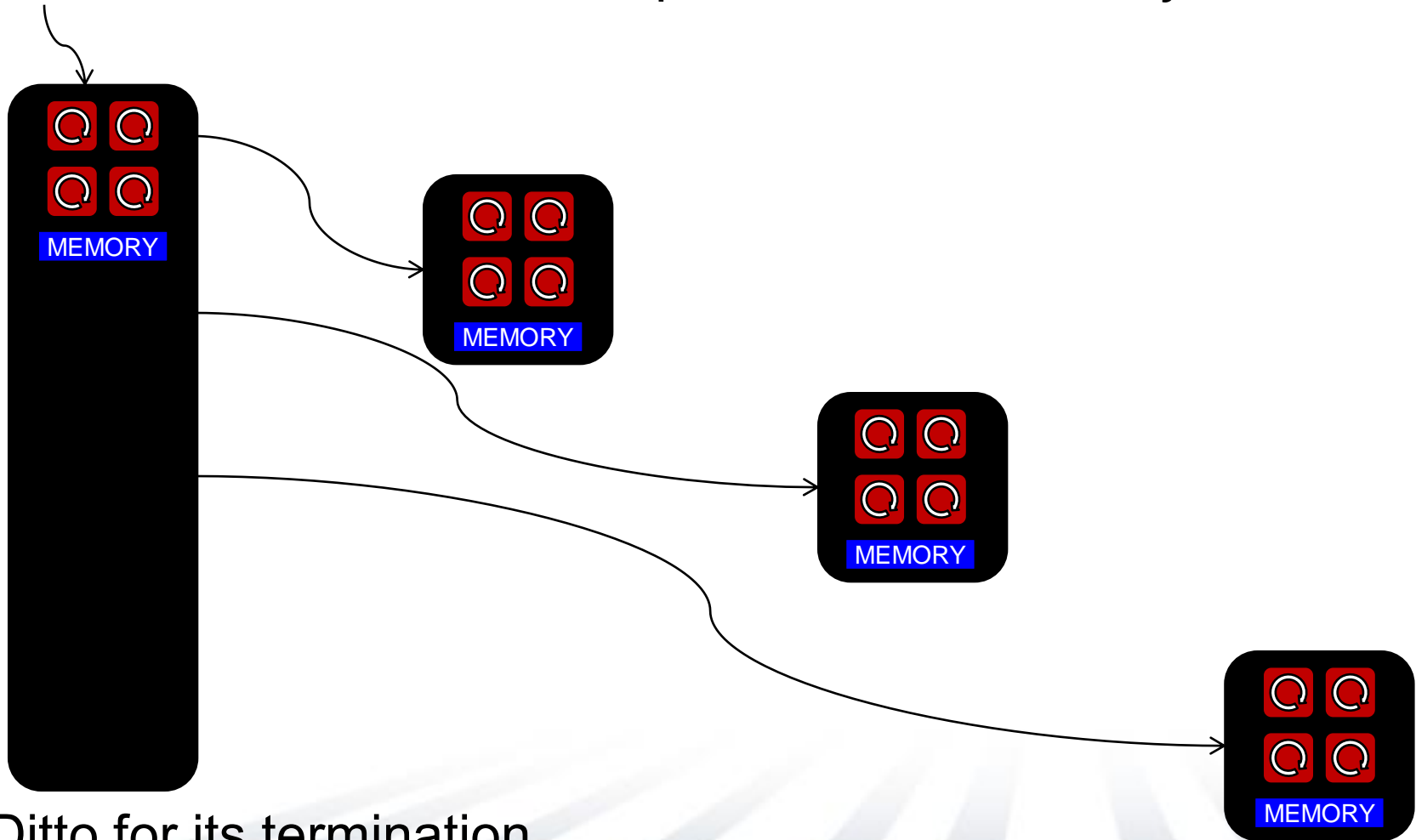
<i>STREAM Triad characteristic</i>	<i>value</i>
per-node problem size	85,985,408
per-node memory required	1.92 GB
percent of available memory	25.0%

# Chapel STREAM Performance



# Why doesn't Chapel scale perfectly?

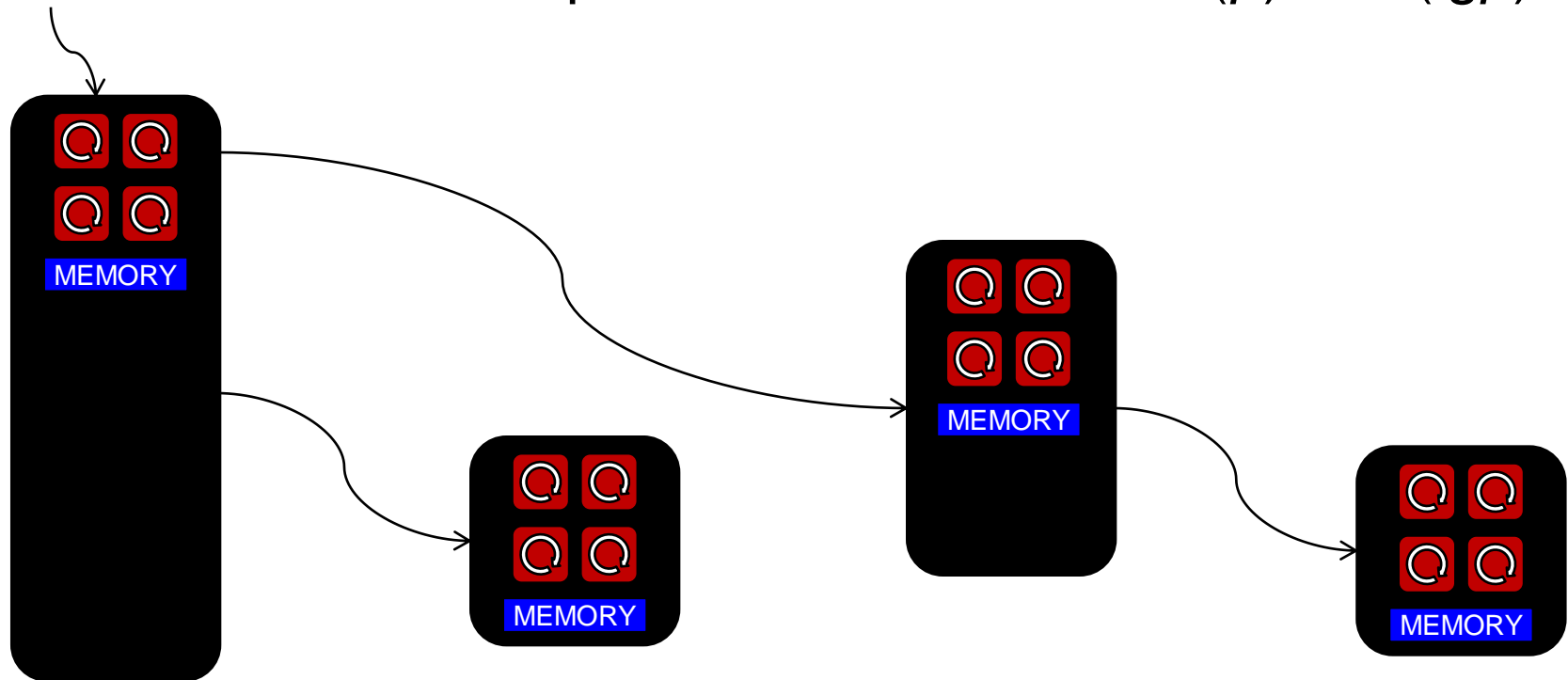
- Because Block1D's current parallel iterator is *very naive*...



- Ditto for its termination...

# Strategies for improvement

- Use tree-based startup/teardown to convert  $O(p)$  to  $O(\lg p)$



- *Or:* Have compiler optimize code to use SPMD exec. model
  - reduces  $O(\lg p)$  to  $O(1)$  by amortizing into program startup/teardown

# SPMD-style Chapel

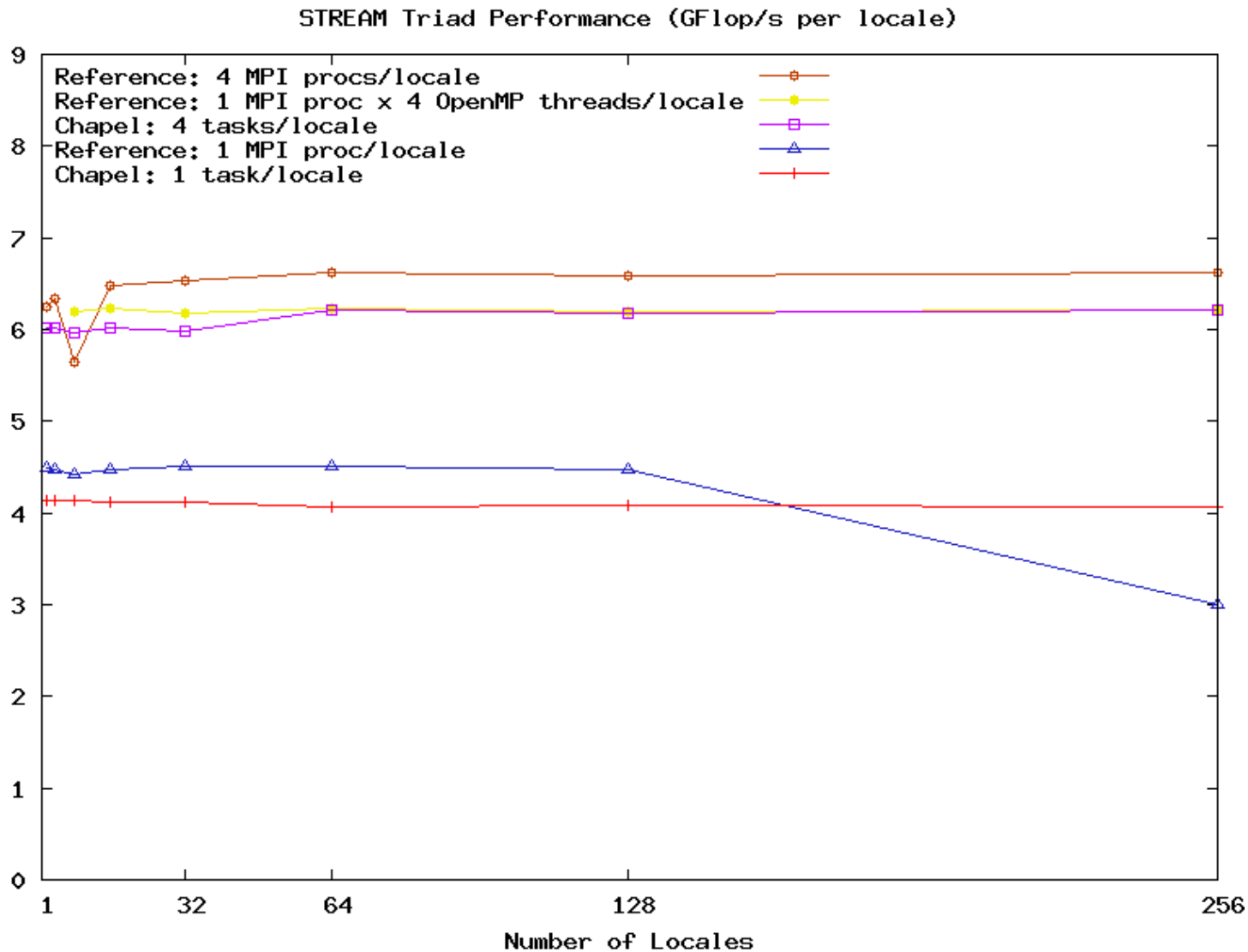
- In the meantime, users can code in SPMD like the MPI version using Chapel's support for *multiresolution programming*:

```
var localGBs: [LocaleSpace] real;

coforall loc in Locales do
on loc {
const myProblemSpace: domain(1, int(64))
    = BlockPartition(ProblemSpace, here.id, numLocales);
var myA, myB, myC: [myProblemSpace] real(64);
const startTime = getCurrentTime();
local {
for (a, b, c) in (myA, myB, myC) do
    a = b + alpha * c;
}
const execTime = getCurrentTime() - startTime;
localGBs(here.id) = timeToGFlops(execTime);
}

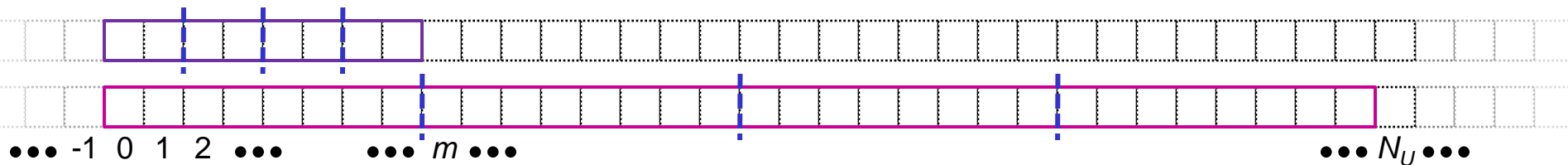
const avgGBs = (+ reduce localGBs) / numLocales;
```

# SPMD Chapel Performance

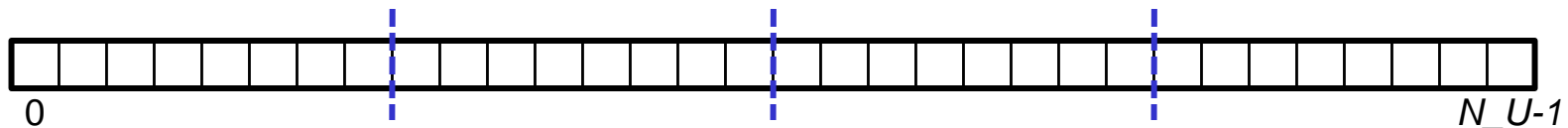
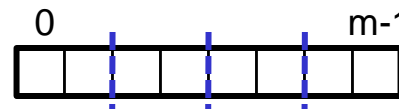


# RA Declarations in Chapel

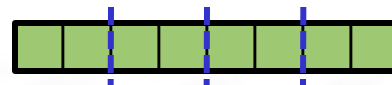
```
const TableDist = new Block1D(bbox=[0..m-1], tasksPerLocale=...),
UpdateDist = new Block1D(bbox=[0..N_U-1], tasksPerLocale=...);
```



```
const TableSpace: domain(1, uint(64)) distributed TableDist = [0..m-1],
Updates: domain(1, uint(64)) distributed UpdateDist = [0..N_U-1];
```



```
var T: [TableSpace] uint(64);
```

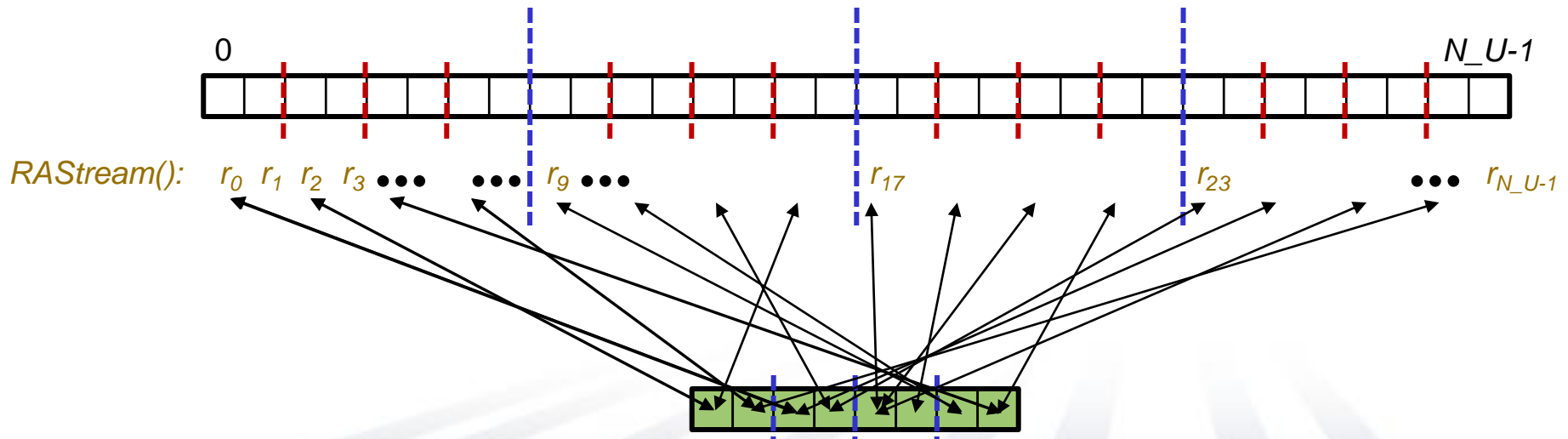


# RA Computation in Chapel

```
const TableSpace: domain(1, uint(64)) distributedTableDist = [0..m-1],
      Updates: domain(1, uint(64)) distributedUpdateDist = [0..N_U-1];
```

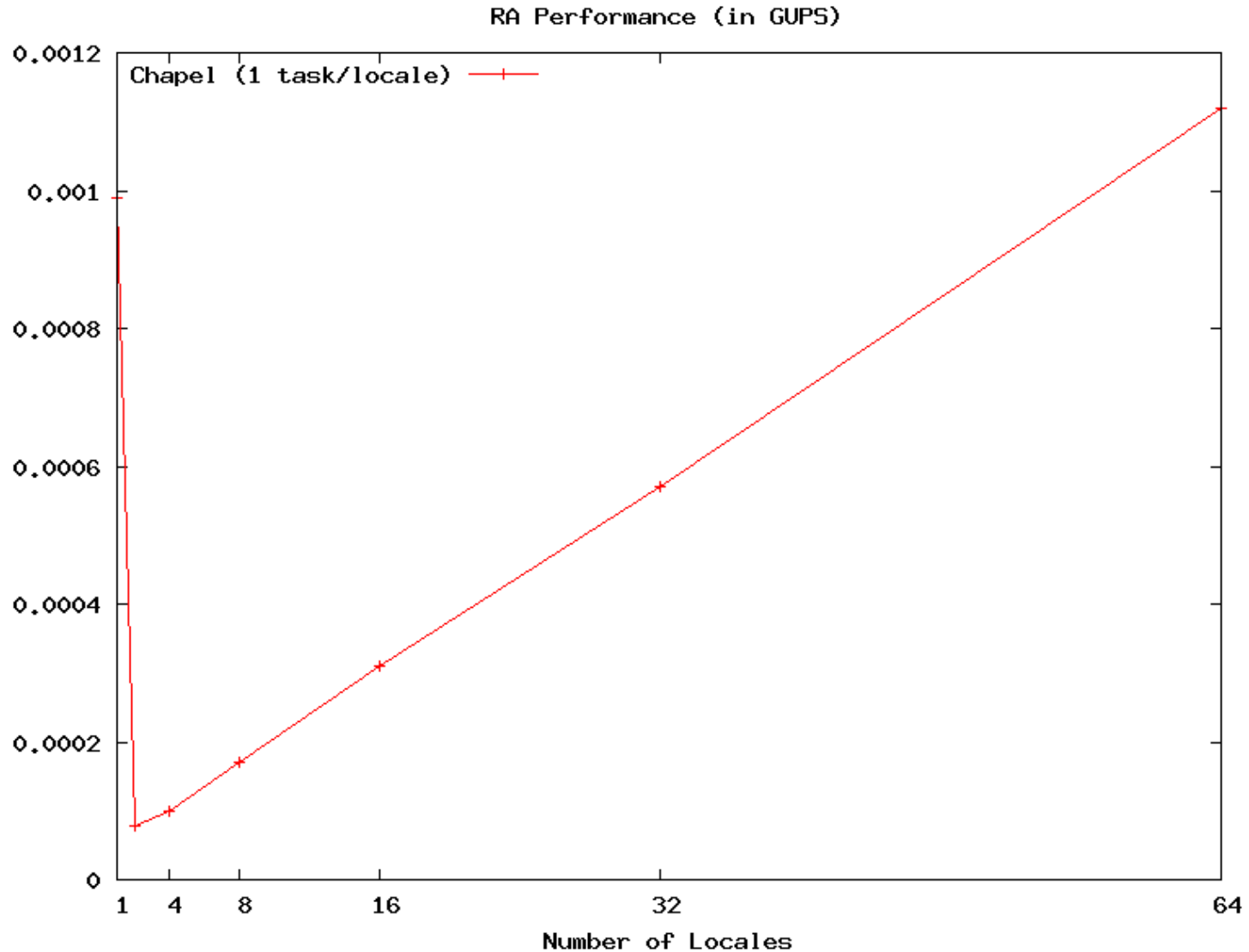
```
var T: [TableSpace] uint(64);
```

```
forall (_, r) in (Updates, RASTream()) do
  on T(r&indexMask) do
    T(r&indexMask) ^= r;
```





# RA Performance in Chapel



# FFT and HPL Status

- FFT :
  - not yet running on distributed memory
    - Block1D not yet rich enough to support slicing, re-indexing
  - have made a big effort to reclaim descriptor memory from slicing
    - can now run full problem size
  
- HPL:
  - not yet running on distributed memory
    - need to add block-cyclic, dimensional, and replicated distributions
  - current version written to be locality-aware
  
- All four of these codes are very clean and should serve as great references to others attempting the HPC Challenge

# Acknowledgements

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*Thanks also to our many colleagues who have helped us reach this point!*

# Summary

- Chapel is scaling on dist. memory machines, if not perfectly
  - more importantly, scalability limiters are known and addressable
- Chapel achieved its first TeraFlop
- Chapel has started to demonstrate user-defined distributions
  - Recall that these have only been working for two months
  - (and a busy two months at that: first public release, two tutorials, ...)
- See you at HPCC 2009!

In the meantime, download Chapel, try it out,  
and please give us your feedback:

<http://chapel.cs.washington.edu>

(our HPCC codes and report are available within the release)