

# Chapel

## Cray Cascade's High Productivity Language

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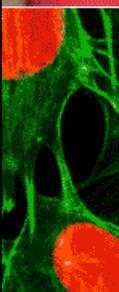
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This Presentation May Contain Some Preliminary Information, Subject To Change



# Chapel Contributors

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# Chapel's Context

**HPCS** = High *Productivity* Computing Systems  
(a DARPA program)

**Overall Goal:** Increase productivity by 10× by 2010

**Productivity** = Programmability  
+ Performance  
+ Portability  
+ Robustness

**Result must be...**

- ...revolutionary, not evolutionary
- ...marketable product

**Phase II Competitors (7/03-7/06):** Cray (Cascade), IBM, Sun

# Chapel Design Objectives

- a *global* view of computation
- support for general parallelism
  - data- and task-parallel; nested parallelism
- clean separation of algorithm and implementation
- broad-market language features
  - OOP, GC, latent types, overloading, generic functions/types, ...
- data abstractions
  - sparse arrays, hash tables, sets, graphs, ...
- good performance
- portability
- interoperability with existing codes

# Outline

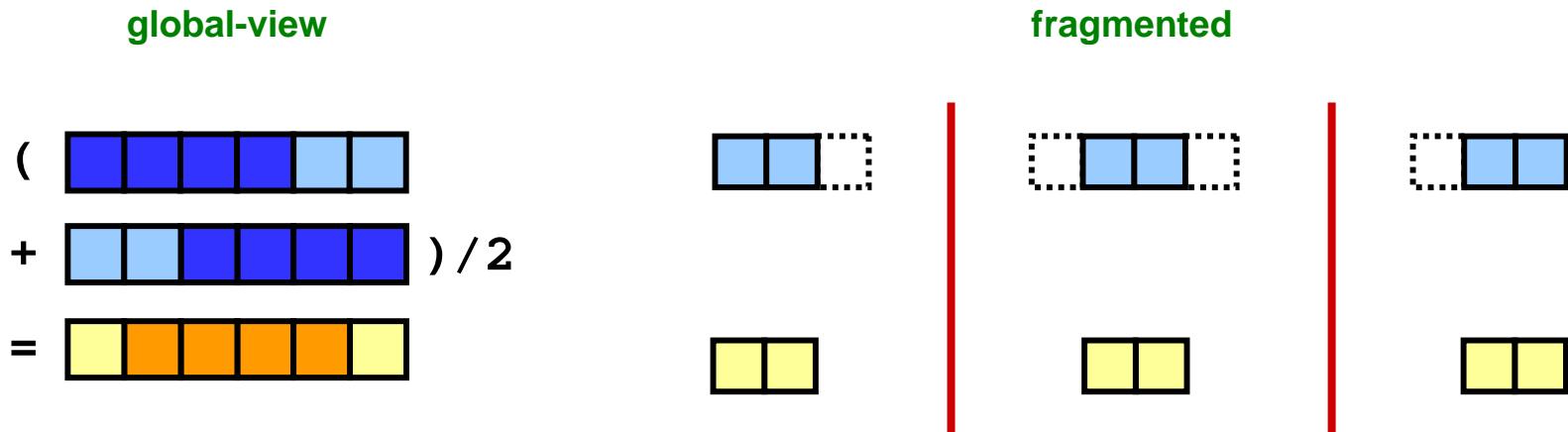
- Chapel Motivation & Foundations
  - ✓ Context and objectives for Chapel
  - Programming models and productivity
- Chapel Overview
- Chapel Activities and Plans

# Parallel Programming Models

- *Fragmented Programming Models:*
  - Programmers *must* program on a task-by-task basis:
    - break distributed data structures into per-task chunks:
    - break work into per-task iterations/control flow
- *Global-view Programming Models:*
  - Programmers need not program task-by-task
    - access distributed data structures as though local
    - introduce parallelism using language keywords
    - burden of decomposition shifts to compiler/runtime
    - user may guide this process via language constructs

# Global-view vs. Fragmented

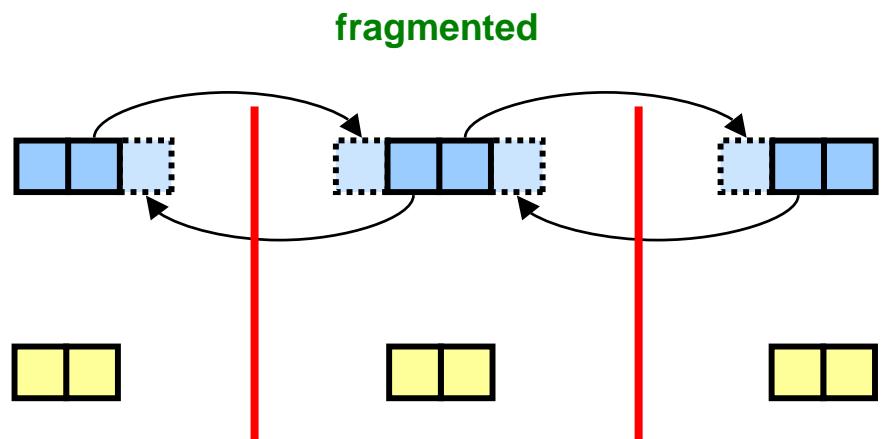
- **Example:** “Apply 3-pt stencil to vector”



# Global-view vs. Fragmented

- **Example:** “Apply 3-pt stencil to vector”

global-view

$$\begin{aligned} & ( \begin{array}{ccccc} \text{blue} & \text{blue} & \text{blue} & \text{blue} & \text{light blue} \\ \hline \end{array} ) \\ & + \begin{array}{ccccc} \text{light blue} & \text{light blue} & \text{blue} & \text{blue} & \text{blue} \end{array} ) / 2 \\ = & \begin{array}{ccccc} \text{yellow} & \text{orange} & \text{orange} & \text{orange} & \text{yellow} \end{array} \end{aligned}$$


# Global-view vs. Fragmented

- **Example:** “Apply 3-pt stencil to vector”

global-view

$$\begin{aligned} & ( \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{light blue}} \boxed{\text{light blue}} ) \\ & + ( \boxed{\text{light blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} ) / 2 \\ & = \boxed{\text{yellow}} \boxed{\text{orange}} \boxed{\text{orange}} \boxed{\text{orange}} \boxed{\text{orange}} \boxed{\text{yellow}} \end{aligned}$$

fragmented

$$\begin{array}{c|c|c|c} \begin{array}{c} ( \boxed{\text{blue}} \boxed{\text{light blue}} \boxed{\text{light blue}} ) \\ + ( \boxed{\text{light blue}} \boxed{\text{light blue}} \boxed{\text{blue}} ) / 2 \\ = \boxed{\text{yellow}} \boxed{\text{orange}} \end{array} & \boxed{\text{red}} & \begin{array}{c} ( \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{light blue}} \boxed{\text{light blue}} ) \\ + ( \boxed{\text{light blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} ) / 2 \\ = \boxed{\text{orange}} \boxed{\text{orange}} \end{array} & \boxed{\text{red}} & \begin{array}{c} ( \boxed{\text{light blue}} \boxed{\text{blue}} \boxed{\text{light blue}} \boxed{\text{light blue}} ) \\ + ( \boxed{\text{light blue}} \boxed{\text{blue}} \boxed{\text{blue}} \boxed{\text{blue}} ) / 2 \\ = \boxed{\text{orange}} \boxed{\text{yellow}} \end{array} \\ \hline \end{array}$$

# Global-view vs. Fragmented

- **Example:** “Apply 3-pt stencil to vector”

## global-view

```
var n: int = 1000;
var a, b: [1..n] float;

forall i in (2..n-1) {
    b(i) = (a(i-1) + a(i+1))/2;
}
```

Assumes *numProcs* divides *n*;  
a more general version would  
require additional effort

## fragmented

```
var n: int = 1000;
var locN: int = n/numProcs;
var a, b: [0..locN+1] float;
var innerLo: int = 1;
var innerHi: int = locN;

if (iHaveRightNeighbor) {
    send(right, a(locN));
    recv(right, a(locN+1));
} else {
    innerHi = locN-1;
}
if (iHaveLeftNeighbor) {
    send(left, a(1));
    recv(left, a(0));
} else {
    innerLo = 2;
}
forall i in (innerLo..innerHi) {
    b(i) = (a(i-1) + a(i+1))/2;
}
```

# Global-view vs. Fragmented

- **Example:** “Apply 3-pt stencil to vector”

## fragmented (pseudocode + MPI)

```
var n: int = 1000, locN: int = n/numProcs;
var a, b: [0..locN+1] float;
var innerLo: int = 1, innerHi: int = locN;
var numProcs, myPE: int;
var retval: int;
var status: MPI_Status;

MPI_Comm_size(MPI_COMM_WORLD, &numProcs);
MPI_Comm_rank(MPI_COMM_WORLD, &myPE);
if (myPE < numProcs-1) {
    retval = MPI_Send(&(a(locN)), 1, MPI_FLOAT, myPE+1, 0, MPI_COMM_WORLD);
    if (retval != MPI_SUCCESS) { handleError(retval); }
    retval = MPI_Recv(&(a(locN+1)), 1, MPI_FLOAT, myPE+1, 1, MPI_COMM_WORLD, &status);
    if (retval != MPI_SUCCESS) { handleErrorWithStatus(retval, status); }
} else
    innerHi = locN-1;
if (myPE > 0) {
    retval = MPI_Send(&(a(1)), 1, MPI_FLOAT, myPE-1, 1, MPI_COMM_WORLD);
    if (retval != MPI_SUCCESS) { handleError(retval); }
    retval = MPI_Recv(&(a(0)), 1, MPI_FLOAT, myPE-1, 0, MPI_COMM_WORLD, &status);
    if (retval != MPI_SUCCESS) { handleErrorWithStatus(retval, status); }
} else
    innerLo = 2;
forall i in (innerLo..innerHi) {
    b(i) = (a(i-1) + a(i+1))/2;
}
```

Communication becomes  
geometrically more complex for  
higher-dimensional arrays

# Fortran+MPI 3D NAS MG Stencil

```

subroutine comm3(u,n1,n2,n3,kk)
use caf_intrinsics
implicit none
include 'cafnpb.h'
include 'globals.h'
integer n1, n2, n3, kk
double precision u(n1,n2,n3)
integer axis
if(.not. dead(kk)) then
do axis = 1, 3
  if( nproc .ne. 1 ) then
    call sync_all()
    call give3( axis, +1, u, n1, n2, n3, kk )
    call sync_all()
    call take3( axis, -1, u, n1, n2, n3 )
    call take3( axis, +1, u, n1, n2, n3 )
  else
    call commmp( axis, u, n1, n2, n3, kk )
  endif
endif
else
do axis = 1, 3
  call sync_all()
  call sync_all()
enddo
call zero3(u,n1,n2,n3)
endif
return
end

subroutine give3( axis, dir, u, n1, n2, n3, k )
use caf_intrinsics
implicit none
include 'cafnpb.h'
include 'globals.h'
integer axis, dir, n1, n2, n3, k, ierr
double precision u( n1, n2, n3 )
integer i3, i2, ii, buff_len,buff_id
buff_id = 2 + dir
buff_len = 0
if( axis .eq. 1 )then
  if( dir .eq. -1 )then
    do i3=2,n1
      do i2=2,n2
        buff_len = buff_len + 1
        buff(buff_len, buff_id ) = u( 2, i2,i3 )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  else if( dir .eq. +1 ) then
    do i3=2,n1
      do i2=2,n2
        buff_len = buff_len + 1
        buff(buff_len, buff_id ) = u( i1,n2-1,i3 )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  endif
endif
if( axis .eq. 2 )then
  if( dir .eq. -1)then
    do i3=2,n1
      do i2=2,n2
        indx = indx + 1
        u(i1,i2,i3) = buff(indx, buff_id )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  else if( dir .eq. +1 ) then
    do i3=2,n1
      do i2=2,n2
        indx = indx + 1
        u(i1,i2,i3) = buff(indx, buff_id )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  endif
endif
if( axis .eq. 3 )then
  if( dir .eq. -1)then
    do i3=2,n1
      do i2=2,n2
        indx = indx + 1
        u(i1,i2,i3) = buff(indx, buff_id )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  else if( dir .eq. +1 ) then
    do i3=2,n1
      do i2=2,n2
        indx = indx + 1
        u(i1,i2,i3) = buff(indx, buff_id )
      enddo
    enddo
    buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
    buff(1:buff_len,buff_id)
  endif
endif
return
end

subroutine give3( axis, dir, u, n1, n2, n3, k )
use caf_intrinsics
implicit none
include 'cafnpb.h'
include 'globals.h'
integer axis, dir, n1, n2, n3, k, ierr
double precision u( n1, n2, n3 )
integer i3, i2, ii, buff_len,buff_id
buff_id = 2 + dir
buff_len = 0
if( axis .eq. 1 )then
  if( dir .eq. -1 )then
    do i3=2,n1
      do i2=2,n2
        buff_len = buff_len + 1
        double precision u( n1, n2, n3 )
      enddo
    enddo
    integer i3, i2, ii, buff_len,buff_id
    integer i1, kk, indx
    dir = -1
    buff_id = 3 + dir
    buff_len = mm2
    do i1=mm2
      buff(i1,buff_id) = 0.0D0
    enddo
    integer i3, i2, ii
    dir = +1
    buff_id = 3 + dir
    buff_len = mm2
    do i1=1,nm2
      buff(i1,buff_id) = 0.0D0
    enddo
    dir = +1
    buff_id = 2 + dir
    buff_len = 0
    if( axis .eq. 1 )then
      do i3=2,n1
        do i2=2,n2
          indx = indx + 1
          u(i1,i2,i3) = buff(indx, buff_id )
        enddo
      enddo
      buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
      buff(1:buff_len,buff_id)
    else if( dir .eq. +1 ) then
      do i3=2,n1
        do i2=2,n2
          indx = indx + 1
          u(i1,i2,i3) = buff(indx, buff_id )
        enddo
      enddo
      buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
      buff(1:buff_len,buff_id)
    endif
    if( axis .eq. 2 )then
      if( dir .eq. -1)then
        do i3=2,n1
          do i2=2,n2
            indx = indx + 1
            u(i1,i2,i3) = buff(indx, buff_id )
          enddo
        enddo
        buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
        buff(1:buff_len,buff_id)
      else if( dir .eq. +1 ) then
        do i3=2,n1
          do i2=2,n2
            indx = indx + 1
            u(i1,i2,i3) = buff(indx, buff_id )
          enddo
        enddo
        buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
        buff(1:buff_len,buff_id)
      endif
    endif
    if( axis .eq. 3 )then
      if( dir .eq. -1)then
        do i3=2,n1
          do i2=2,n2
            indx = indx + 1
            u(i1,i2,i3) = buff(indx, buff_id )
          enddo
        enddo
        buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
        buff(1:buff_len,buff_id)
      else if( dir .eq. +1 ) then
        do i3=2,n1
          do i2=2,n2
            indx = indx + 1
            u(i1,i2,i3) = buff(indx, buff_id )
          enddo
        enddo
        buff(1:buff_len,buff_id+1)[nbr(axis,dir,k)] =
        buff(1:buff_len,buff_id)
      endif
    endif
    return
  end

subroutine commmp( axis, u, n1, n2, n3, kk )
use caf_intrinsics
implicit none
include 'cafnpb.h'
include 'globals.h'
integer m1k, m2k, m3k, s,m1j,m2j,m3j,jk
double precision r(m1k,m2k,m3k), s(m1j,m2j,m3j)
integer j3, 32, jl, i3, i2, il, d1, d2, d3, j
double precision xl(m), yl(m), x2,y2
if(mlk.eq.3)then
  d1 = 2
else
  d1 = 1
endif
if(m2k.eq.3)then
  d2 = 2
else
  d2 = 1
endif
if(m3k.eq.3)then
  d3 = 2
else
  d3 = 1
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      y1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(j1,j2,j3) =
      > + 0.35D0 * ( r(i1-1,i2,i3) + r(i1+1,i2,i3) + x2 )
      > + 0.125D0 * ( x1(i1-1) + x1(i1+1) + y2 )
      > + 0.0625D0 * ( y1(i1-1) + y1(i1+1) )
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(j1,j2,j3) =
      > + 0.35D0 * ( r(i1-1,i2,i3) + r(i1+1,i2,i3) + x2 )
      > + 0.125D0 * ( x1(i1-1) + x1(i1+1) + y2 )
      > + 0.0625D0 * ( y1(i1-1) + y1(i1+1) )
    enddo
  enddo
endif
end
j = 1
call comm3(s,m1j,m2j,m3j,j)
return
end

subroutine rpxj3(r,m1k,m2k,m3k,s,m1j,m2j,m3j,k)
implicit none
include 'cafnpb.h'
include 'globals.h'
integer mlk, m2k, m3k, s,m1j,m2j,m3j,jk
double precision r(m1k,m2k,m3k), s(m1j,m2j,m3j)
integer j3, 32, jl, i3, i2, il, d1, d2, d3, j
double precision xl(m), yl(m), x2,y2
if(mlk.eq.3)then
  d1 = 2
else
  d1 = 1
endif
if(m2k.eq.3)then
  d2 = 2
else
  d2 = 1
endif
if(m3k.eq.3)then
  d3 = 2
else
  d3 = 1
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      y1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(j1,j2,j3) =
      > + 0.35D0 * ( r(i1-1,i2,i3) + r(i1+1,i2,i3) + x2 )
      > + 0.125D0 * ( x1(i1-1) + x1(i1+1) + y2 )
      > + 0.0625D0 * ( y1(i1-1) + y1(i1+1) )
    enddo
  enddo
endif
do j3=2,m3j-1
  i3 = i3+j3*d3
  do j2=2,m2j-1
    i2 = 2*j2-d2
    do j1=2,m1j
      i1 = 2*j1-d1
      x1(i1-1) = r(i1-1,i2-1,i3-1) + r(i1-1,i2+1,i3 )
      > + r(i1-1,i2-1,i3+1) + r(i1-1,i2+1,i3+1)
      x2 = r(i1, i2-1,i3 ) + r(i1, i2+1,i3 )
      > + r(i1, i2-1,i3-1) + r(i1, i2+1,i3-1)
      y1(j1,j2,j3) =
      > + 0.35D0 * ( r(i1-1,i2,i3) + r(i1+1,i2,i3) + x2 )
      > + 0.125D0 * ( x1(i1-1) + x1(i1+1) + y2 )
      > + 0.0625D0 * ( y1(i1-1) + y1(i1+1) )
    enddo
  enddo
endif
end
j = 1
call comm3(s,m1j,m2j,m3j,j)
return
end

```

# Chapel 3D NAS MG Stencil

```
param coeff: domain(1) = [0..3]; // for 4 unique weight values
param Stencil: domain(3) = [-1..1, -1..1, -1..1]; // 27-points

function rprj3(S, R) {
    param w: [coeff] float = (/0.5, 0.25, 0.125, 0.0625/);
    param w3d: [(i,j,k) in Stencil] float
        = w((i!=0) + (j!=0) + (k!=0));
    const SD = S.domain,
          Rstr = R.stride;

    S = [ijk in SD] sum reduce [off in Stencil]
        (w3d(off) * R(ijk + Rstr*off));
}
```

# Fragmented Language Summary

- Fragmented programming models...
  - ...manage per-task details in-line with the computation
    - per-task local bounds, data structures
    - communication, synchronization
  - ...are our main parallel programmability limiter today

# Fragmented Language Summary

- Fragmented programming models...
  - ...tend to be easier to compile than global-view languages
    - at minimum, only need a good node compiler
  - ...deserve credit for the majority of the community's parallel application successes to date

# Global-View Language Summary

- Single-processor languages are trivially global-view
  - Matlab, Java, Python, Perl, C, C++, Fortran, ...
- Parallel global-view languages have been developed...
  - HPF (High Performance Fortran), ZPL, Sisal, NESL, Cilk, Cray MTA extensions to C/Fortran, ...
- ...yet most have not achieved widespread adoption
  - reasons why are as varied as the languages themselves
- Chapel has been designed...
  - ...to support global-view programming
  - ...with experience from preceding global-view languages

# Outline

- ✓ Chapel Motivation & Foundations
- Chapel Overview
- Chapel Activities and Plans

# What is Chapel?

- *Chapel*: Cascade High-Productivity Language
- Overall goal: “Solve the parallel programming problem”
  - simplify the creation of parallel programs
  - support their evolution to extreme-performance, production-grade codes
  - emphasize generality
- Motivating Language Technologies:
  - global-view multithreaded parallel programming
  - locality-aware programming

# Multithreaded Parallel Programming

- Virtualization of threads
  - *i.e.*, no fork/join, naming of threads
- Abstractions for data and task parallelism
  - *data*: domains, arrays, iterators, ...
  - *task*: cobegins, atomic transactions, sync variables, ...
- Composition of parallelism
- Global view of computation, data structures

# Data Parallelism: Domains

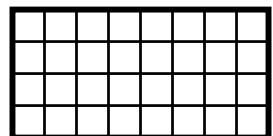
- *domain*: an index set
  - specifies size and shape of arrays
  - supports sequential and parallel iteration
  - potentially decomposed across locales
- Three main classes:
  - *arithmetic*: indices are Cartesian tuples
    - rectilinear, multidimensional, optionally strided and/or sparse
  - *indefinite*: indices serve as hash keys
    - supports hash tables, associative arrays, dictionaries
  - *opaque*: indices are anonymous
    - supports sets, graph-based computations
- Chapel's fundamental concept for data parallelism

# Simple Domain Declarations

```
var m: int = 4;
```

```
var n: int = 8;
```

```
var D: domain(2) = [1..m, 1..n];
```



$D$

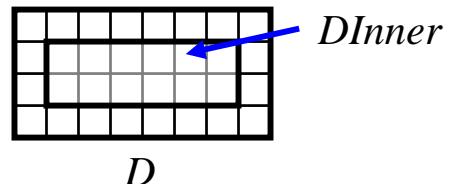
# Simple Domain Declarations

```
var m: int = 4;
```

```
var n: int = 8;
```

```
var D: domain(2) = [1..m, 1..n];
```

```
var DInner: subdomain(D) = [2..m-1, 2..n-1];
```



# Domain Uses

- Declaring arrays:

```
var A, B: [D] float;
```

- Sub-array references:

```
A(DInner) = B(DInner);
```

- Iteration:

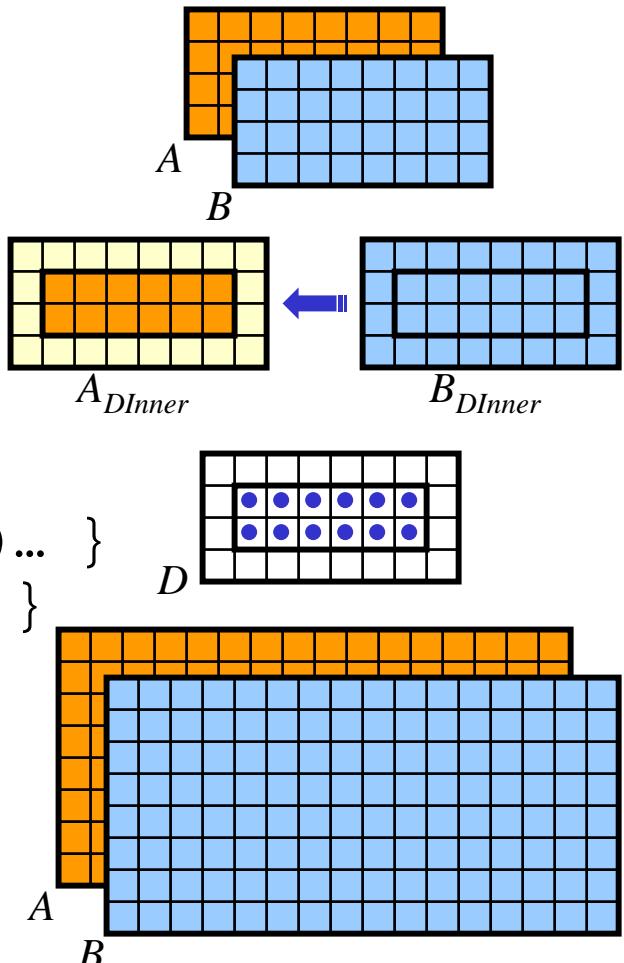
```
forall (i,j) in DInner { ...A(i,j)... }
```

or: **forall** ind in DInner { ...A(ind)... }

or: [ind **in** DInner] ...A(ind)...

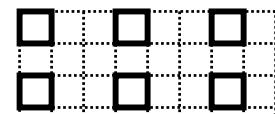
- Array reallocation:

```
D = [1..2*m, 1..2*n];
```



# Other Arithmetic Domains

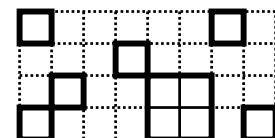
```
var StridedD: subdomain(D) = D by (2,3);
```



*StridedD*

```
var indexList: seq(index(D)) = ...;
```

```
var SparseD: sparse subdomain(D) = indexList;
```



*SparseD*

# Task Parallelism

- *co-begins*: indicate statements that may run in parallel:

```
computePivot(lo, hi, data);  
cobegin {  
    Quicksort(lo, pivot, data);  
    Quicksort(pivot, hi, data);  
}
```

```
cobegin {  
    ComputeTaskA(...);  
    ComputeTaskB(...);  
}
```

- *atomic sections*: support atomic transactions

```
atomic {  
    newnode.next = insertpt;  
    newnode.prev = insertpt.prev;  
    insertpt.prev.next = newnode;  
    insertpt.prev = newnode;  
}
```

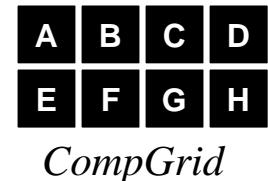
- *sync and single-assignment variables*: synchronize tasks
  - similar to Cray MTA C/Fortran

# Locality-aware Programming

- *locale*: architectural unit of storage and processing
- programmer specifies number of locales on executable command-line

```
prompt> myChapelProg -nl=8
```
- Chapel programs are provided with built-in locale array:  
`const Locales: [1..numLocales] locale;`
- Users may use it to create their own locale arrays:

```
var CompGrid: [1..GridRows, 1..GridCols] locale = ...;
```



```
var TaskALocs: [1..numTaskALocs] locale = Locales(1..2);
```

```
var TaskBLocs: [1..numTaskBLocs] locale = Locales(3..numLocales);
```



*TaskALocs*

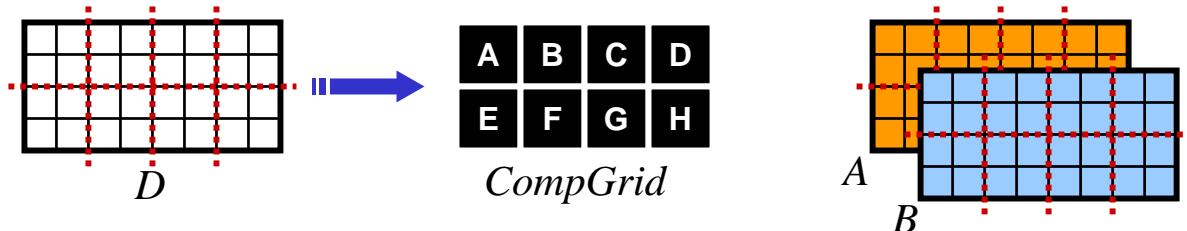


*TaskBLocs*

# Data Distribution

- domains may be distributed across locales

```
var D: domain(2) distributed(Block(2) on CompGrid) = ...;
```



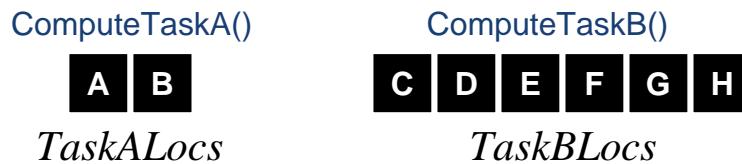
- Distributions specify...
  - ...mapping of indices to locales
  - ...per-locale storage layout of domain indices and array elements
- Distributions implemented as a class hierarchy
  - Chapel provides a number of standard distributions
  - Users may also write their own

one of our biggest challenges

# Computation Distribution

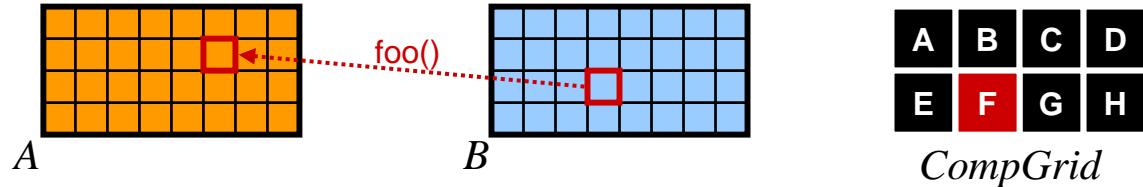
- “on” keyword binds computation to locale(s):

```
cobegin {  
    on TaskALocs do ComputeTaskA(...);  
    on TaskBLocs do ComputeTaskB(...);  
}
```



- “on” can also be used in a data-driven manner:

```
forall (i,j) in D {  
    on B(j/2,i*2) do A(i,j) = foo(B(j/2,i*2));  
}
```



# Chapel Challenges

- User Acceptance
  - True of any new language
  - Skeptical audience
- Commodity Architecture Implementation
  - Chapel designed with idealized architecture in mind
  - Clusters are not ideal in many respects
  - Results in implementation and performance challenges
- And many others as well...

# Outline

- ✓ Chapel Motivation & Foundations
- ✓ Chapel Overview
- Chapel Activities and Plans

# Phase II Activities

- 2003-2006:
  - Application studies to drive language design
    - HPCC, NPB, SSCA benchmarks
    - kernels from Cray customer applications
    - other kernels of interest (connected components, FMM)
  - Design and specification of Chapel language
  - Implementation work on portable Chapel prototype
  - Outreach to inform users and get feedback
    - government: LANL, Sandia, LLNL, ORNL, JPL, NITRD
    - conferences: ICS, PPoPP, LCPC, PGAS, HIPS, HPL, LaR
    - mainstream industry: Microsoft (w/ AMD attendance)
    - HPCS: biannual reviews, SW productivity meetings

# What's next?

- HPCS phase III
  - July 2006 – December 2010
  - 2 vendors expected to be funded
  - proposals submitted May 5th
- HPCS Language Effort forking off
  - all 3 phase II language teams eligible for phase III
  - High Productivity Language Systems (HPLS) team
    - language experts/enthusiasts from national labs, academia
    - to study, evaluate the vendor languages, report to DARPA
    - July 2006 – December 2007
  - DARPA hopes...
    - ...that a language consortium will emerge from this effort
    - ...to involve mainstream computing vendors as well
    - ...to avoid repeating mistakes of the past (Ada, HPF, ...)

# Proposed Phase III Activities

- Short-term (2006-2007):
  - support user evaluations of Chapel
    - HPCS mission partners
    - HPLS language evaluation team
    - software productivity team
    - other potential user communities
  - continue Chapel implementation
  - capture application studies as tutorials
  - revise language as suggested by these activities
- Longer-term (2008-2010):
  - participate in HPLS consortium language efforts
  - help build support for language in community
  - fold HPLS language into Cascade compiler, tools

# Summary

- Chapel is being designed to...
  - ...enhance programmer productivity
  - ...address a wide range of HEC algorithms
- Via high-level, extensible abstractions for...
  - ...multithreaded parallel programming
  - ...locality-aware programming
- Status:
  - *draft* language specification available at:  
<http://chapel.cs.washington.edu>
  - Open source implementation proceeding apace
  - Your feedback desired!