

ONE-DAY CHAPEL TUTORIAL SESSION 2: CHAPEL BASICS

Chapel Team October 16, 2023

ONE DAY CHAPEL TUTORIAL

- 9-10:30: Getting started using Chapel for parallel programming
- 10:30-10:45: break
- 10:45-12:15: Chapel basics in the context of the n-body example code
- 12:15-1:15: lunch
- 1:15-2:45: Distributed and shared-memory parallelism especially w/arrays (data parallelism)
- 2:45-3:00: break
- 3:00-4:30: More parallelism including for asynchronous parallelism (task parallelism)
- 4:30-5:00: Wrap-up including gathering further questions from attendees

OUTLINE: CHAPEL BASICS

- Running Example: n-body computation (Hands On)
- Variables, Constants, and Operators
- Records and Classes
- Tuples
- Arrays
- Writing out Tuples, Records, and Arrays (Hands On)
- Main() Procedure
- Ranges and basic control flow
- Procedures and iterators
- Where might we parallelize the n-body computation? (Hands On)



N-BODY IN CHAPEL (WHERE N == 5)

- A serial computation
- From the Computer Language Benchmarks Game
 - Chapel implementation in release under examples/benchmarks/shootout/nbody.chpl
- Computes the influence of 5 bodies on one another
 - The Sun, Jupiter, Saturn, Uranus, Neptune
- Executes for a user-specifiable number of timesteps

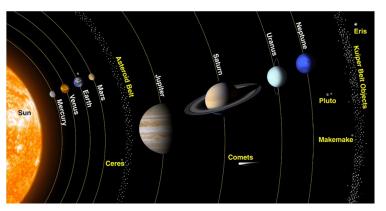


Image source: http://spaceplace.nasa.gov/review/ice-dwarf/solar-system-lrg.png

HANDS ON: COMPILING AND RUNNING N-BODY



Things to try

```
chpl nbody.chpl
time ./nbody -nl 1
time ./nbody -nl 1 -n=100000

chpl --fast nbody.chpl
time ./nbody -nl 1
time ./nbody -nl 1
time ./nbody -nl 1 -n=100000
```

```
// number of timesteps to simulate
config const n = 10000;
...
```

Key concepts

- *nix 'time' command is an easy way to see how long a program takes to run
- Compile with '--fast' to have 'chpl' compiler generate faster code

VARIABLES, CONSTANTS, AND OPERATORS

5-BODY IN CHAPEL: VARIABLE AND RECORD DECLARATIONS



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
                                           Variable declarations
      daysPerYear = 365.24;
config const numsteps = 10000;
record body {
 var pos: 3*real;
 var v: 3*real;
 var mass: real;
```

VARIABLES, CONSTANTS, AND PARAMETERS

Basic syntax

```
declaration:
  var  identifier [: type] [= init-expr];
  const identifier [: type] [= init-expr];
  param identifier [: type] [= init-expr];
```

Examples

Meaning

- var/const: execution-time variable/constant
- param: compile-time constant
- No init-expr ⇒ initial value is the type's default
- No type ⇒ type is taken from init-expr

PRIMITIVE TYPES

Syntax

Туре	Description	Default Value	Currently-Supported Bit Widths	Default Bit Width
bool	logical value	false		impl. dep.
int	signed integer	0	8, 16, 32, 64	64
uint	unsigned integer	0	8, 16, 32, 64	64
real	real floating point	0.0	32, 64	64
imag	imaginary floating point	0.0i	32, 64	64
complex	complex floating points	0.0 + 0.0i	64, 128	128
string	character string	Ш	N/A	N/A

Examples

```
primitive-type:
   type-name [( bit-width )]
```

CHAPEL'S STATIC TYPE INFERENCE

```
const pi = 3.14,
                 // pi is a real
     coord = 1.2 + 3.4i, // coord is a complex...
      coord2 = pi*coord, //...as is coord2
     proc addem (x, y) { // addem() has generic arguments
  return x + y; // and an inferred return type
var sum = addem(1, pi),
                        // sum is a real
   fullname = addem(name, "ford"); // fullname is a string
writeln((sum, fullname));
```

(4.14, bradford)



nbody.chpl

BASIC OPERATORS AND PRECEDENCE

Operator	Description	Associativity	Overloadable
:	cast	left	yes
**	exponentiation	right	yes
!~	logical and bitwise negation	right	yes
* / %	multiplication, division and modulus	left	yes
(unary) + -	positive identity and negation	right	yes
<< >>	shift left and shift right	left	yes
&	bitwise/logical and	left	yes
^	bitwise/logical xor	left	yes
I	bitwise/logical or	left	yes
+-	addition and subtraction	left	yes
<= >= < >	ordered comparison	left	yes
== !=	equality comparison	left	yes
&&	short-circuiting logical and	left	via isTrue
II	short-circuiting logical or	left	via isTrue



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
                                           Variable declarations
      daysPerYear = 365.24;
config const numsteps = 10000;
record body {
 var pos: 3*real;
 var v: 3*real;
 var mass: real;
```

```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
      daysPerYear = 365.24;
config const numsteps = 10000;
                                          Configuration Variable
record body {
 var pos: 3*real;
 var v: 3*real;
 var mass: real;
```



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
      daysPerYear = 365.24;
config const numsteps = 10000;
                                          Configuration Variable
record body {
 var pos: 3*real;
                                  $ ./nbody --numsteps=100
 var v: 3*real;
 var mass: real;
```

CONFIGS



```
param intSize = 32;
type elementType = real(32);
const epsilon = 0.01:elementType;
var start = 1:int(intSize);
```

CONFIGS

```
config param intSize = 32;
config type elementType = real(32);
config const epsilon = 0.01:elementType;
config var start = 1:int(intSize);
```

```
$ chpl 02-configs.chpl -sintSize=64 -selementType=real
$ ./02-configs-start=2 -nl 1 --epsilon=0.00001
```

```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
      daysPerYear = 365.24;
config const numsteps = 10000;
                                          Configuration Variable
record body {
 var pos: 3*real;
 var v: 3*real;
 var mass: real;
```



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
      daysPerYear = 365.24;
config const numsteps = 10000;
record body { ←
 var pos: 3*real;
                                 Record declaration
 var v: 3*real;
 var mass: real;
```

RECORDS AND CLASSES

RECORDS AND CLASSES

Chapel's object types

- Contain variable definitions (fields)
- Contain procedure & iterator definitions (methods)
- Records: value-based (e.g., assignment copies fields)
- Classes: reference-based (e.g., assignment aliases object)

Example

```
use Math;
record circle {
  var radius: real;
  proc area() {
    return pi*radius**2;
  }
}
```

```
var c1 = new circle(radius=1.0);
var c2 = c1;  //copies c1
c1.radius = 5.0;
writeln(c2.radius);  // prints 1.0
```

RECORDS AND CLASSES



Chapel's object types

- Contain variable definitions (fields)
- Contain procedure & iterator definitions (methods)
- Records: value-based (e.g., assignment copies fields)
- Classes: reference-based (e.g., assignment aliases object)

Example

```
use Math;
class Circle {
  var radius: real;
  proc area() {
    return pi*radius**2;
  }
}
```

CLASSES VS. RECORDS

Classes

- heap-allocated
 - Variables point to objects
 - Support mem. mgmt. policies
- 'reference' semantics
 - compiler will only copy pointers
- support inheritance
- support dynamic dispatch
- identity matters most
- similar to Java classes

Records

- allocated in-place
 - Variables are the objects
 - Always freed at end of scope
- 'value' semantics
 - compiler may introduce copies
- no inheritance
- no dynamic dispatch
- value matters most
- similar to C++ structs
 - (sans pointers)



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
      daysPerYear = 365.24;
config const numsteps = 10000;
record body {
 var pos: 3*real;
 var v: 3*real;
 var mass: real;
                                Tuple type
```

OUTLINE: CHAPEL BASICS

- Running Example: n-body computation (Hands On)
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- Where might we parallelize the n-body computation? (Hands On)

TUPLES (HANDS ON)

TUPLES

Use

- support lightweight grouping of values
 - e.g., passing/returning multiple procedure arguments at once
 - short vectors
 - multidimensional array indices
- support heterogeneous data types

Examples

```
var coord: (int, int, int) = (1, 2, 3);
var coordCopy: 3*int = coord;
var (i1, i2, i3) = coord;
var triple: (int, string, real) = (7, "eight", 9.0);
```



```
const pi = 3.141592653589793,
      solarMass = 4 * pi**2,
                                             Variable declarations
      daysPerYear = 365.24;
config const numsteps = 10000;
                                            Configuration Variable
record body {
 var pos: 3*real;
                                   Record declaration
 var v: 3*real;
 var mass: real;
                                   Tuple type
```

nbody.chpl

```
var bodies =
    [ /* sun */
       new body(mass = solarMass),
       /* jupiter */
       new body (pos = (4.84143144246472090e+00,
                       -1.16032004402742839e+00,
                       -1.03622044471123109e-01),
                  v = (1.66007664274403694e-03 * daysPerYear,
                        7.69901118419740425e-03 * daysPerYear,
                       -6.90460016972063023e-05 * daysPerYear),
               mass = 9.54791938424326609e-04 * solarMass),
       /* saturn */
       new body(...),
       /* uranus */
       new body (...),
       /* neptune */
       new body (...)
   ];
```

```
var bodies =
    [ /* sun */
                                         Create a record object
       new body(mass = solarMass),
       /* jupiter */
       new body (pos = (4.84143144246472090e+00,
                        -1.16032004402742839e+00,
                        -1.03622044471123109e-01),
                  v = (1.66007664274403694e-03 * daysPerYear,
                         7.69901118419740425e-03 * daysPerYear,
                        -6.90460016972063023e-05 * daysPerYear),
               mass = 9.54791938424326609e-04 * solarMass),
       /* saturn */
       new body (...),
       /* uranus */
       new body (...),
       /* neptune */
       new body (...)
   ];
```

```
var bodies =
    [ /* sun */
       new body(mass = solarMass),
       /* jupiter */
                                                            Tuple values
       new body (pos = (4.84143144246472090e+00,
                        -1.16032004402742839e+00,
                        -1.03622044471123109e-01),
                  v = (1.66007664274403694e-03 * daysPerYéar,
                        7.69901118419740425e-03 * daysPerYear,
                       -6.90460016972063023e-05 * daysPerYear),
               mass = 9.54791938424326609e-04 * solarMass),
       /* saturn */
       new body (...),
       /* uranus */
       new body (...),
       /* neptune */
       new body (...)
   ];
```

nbody.chpl

```
var bodies =
    [ /* sun */
       new body(mass = solarMass),
       /* jupiter */
       new body (pos = (4.84143144246472090e+00,
                       -1.16032004402742839e+00,
                       -1.03622044471123109e-01),
  Array
                  v = (1.66007664274403694e-03 * daysPerYear,
  value
                        7.69901118419740425e-03 * daysPerYear,
                       -6.90460016972063023e-05 * daysPerYear),
               mass = 9.54791938424326609e-04 * solarMass),
       /* saturn */
       new body (...),
       /* uranus */
       new body (...),
       /* neptune */
       new body (...)
```

ARRAYS

ARRAY TYPES

Syntax

```
array-type:
   [ domain-expr ] elt-type
array-value:
   [elt1, elt2, elt3, ... eltn]
```

Meaning

- array-type: stores an element of elt-type for each index
- array-value: represent the array with these values

Examples

More on arrays in data parallelism section later...



```
var bodies =
    [ /* sun */
                                         Create a record object
       new body(mass = solarMass),
       /* jupiter */
                                                             Tuple values
       new body (pos = (4.84143144246472090e+00),
                        -1.16032004402742839e+00,
                        -1.03622044471123109e-01),
  Array
                   v = (1.66007664274403694e-03 * daysPerYear,
  value
                         7.69901118419740425e-03 * daysPerYear,
                        -6.90460016972063023e-05 * daysPerYear),
               mass = 9.54791938424326609e-04 * solarMass),
       /* saturn */
       new body (...),
       /* uranus */
       new body (...),
       /* neptune */
       new body (...)
```

HANDS ON: WRITING TUPLES, RECORDS, AND ARRAYS



Put a 'writeln("bodies = ", bodies);' into program

```
chpl nbody.chpl
./nbody -nl 1
bodies = (pos = (0.0, 0.0, 0.0), vel = (0.0, 0.0, 0.0),
mass = 39.4784) (pos = (4.84143, -1.16032, -0.103622), vel
= (0.606326, 2.81199, -0.0252184), mass = 0.0376937) (pos
= (8.34337, 4.1248, -0.403523), vel = (-1.01077, 1.82566,
0.00841576), mass = 0.0112863) (pos = (12.8944, -15.1112,
[-0.223308), vel = (1.08279, 0.868713, -0.0108326), mass =
0.00172372) (pos = (15.3797, -25.9193, 0.179259), vel =
(0.979091, 0.594699, -0.034756), mass = 0.00203369)
-0.169075164
-0.169016441
```

MAIN() PROCEDURE

```
proc main() {
  initSun();
  writef("%.9r\n", energy());
  for 1..numsteps do
    advance(0.01);
  writef("%.9r\n", energy());
```

```
Procedure Definition
proc main() {
  initSun();
  writef("%.9r\n", energy());
  for 1..numsteps do
    advance(0.01);
  writef("%.9r\n", energy());
```

```
proc main() {
                                   Procedure Call
  initSun();
  writef("%.9r\n", energy());
  for 1..numsteps do
    advance(0.01);
  writef("%.9r\n", energy());
```

```
proc main() {
  initSun();
  writef("%.9r\n", energy());
                                            Formatted I/O
  for 1..numsteps do
    advance(0.01);
 writef("%.9r\n", energy());
```

```
proc main() {
  initSun();
  writef("%.9r\n", energy());
  for 1..numsteps do
    advance (0.01);
  writef "%.9r\n", energy());
                   Range Value
```

RANGES: INTEGER SEQUENCES

RANGE VALUES: INTEGER SEQUENCES

Syntax

```
range-expr:
[low] .. [high]
```

Definition

Regular sequence of integers

low <= high: low, low+1, low+2, ..., high</pre>

low > high: degenerate (an empty range)

low or high unspecified: unbounded in that direction

Examples

```
1..6 // 1, 2, 3, 4, 5, 6
6..1 // empty
3.. // 3, 4, 5, 6, 7, ...
```

RANGE OPERATORS

```
const r = 1..10;
printVals(r);
printVals(r # 3);
printVals(r by 2);
printVals(r by -2);
printVals(r by 2 # 3);
printVals(r # 3 by 2);
printVals(0.. #n);
proc printVals(r) {
  for i in r do
   write(i, " ");
  writeln();
```

```
1 2 3 4 5 6 7 8 9 10
1 2 3
1 3 5 7 9
10 8 6 4 2
1 3 5
1 3
0 1 2 3 4 ... n-1
```

```
proc main() {
  initSun();
  writef("%.9r\n", energy());
  for 1...numsteps do
    advance(0.01);
  writef("%.9r\n", energy());
                                             Serial for loop
```

BASIC SERIAL CONTROL FLOW

FOR LOOPS

Syntax

```
for-loop:
   for [index-expr in] iteratable-expr { stmt-list }
```

Meaning

- Executes loop body serially, once per loop iteration
- Declares new variables for identifiers in *index-expr*
 - -type and const-ness determined by *iteratable-expr*
 - -iteratable-expr could be a range, array, iterator, iterable object, ...

Examples

```
var A: [1..3] string = [" DO", " RE", " MI"];
for i in 1..3 { write(A[i]); } //DOREMI
for a in A { a += "LA"; } write(A); //DOLA RELA MILA
```

CONTROL FLOW: OTHER FORMS

Conditional statements

```
if cond { computeA(); } else { computeB(); }
```

While loops

```
while cond {
  compute();
}
```

For loops

```
for indices in iteratable-expr {
  compute();
}
```

```
select key {
  when value1 { compute1(); }
  when value2 { compute2(); }
  otherwise { compute3(); }
}
```

CONTROL FLOW: BRACES VS. KEYWORDS

Control flow statements specify bodies using curly brackets (compound statements)

Conditional statements

```
if cond { computeA(); } else { computeB(); }
```

While loops

```
while cond {
  compute();
}
```

For loops

```
for indices in iteratable-expr {
  compute();
}
```

```
select key {
  when value1 { compute1(); }
  when value2 { compute2(); }
  otherwise { compute3(); }
}
```

CONTROL FLOW: BRACES VS. KEYWORDS

They also support keyword-based forms for single-statement cases

Conditional statements

```
if cond then computeA(); else computeB();
```

While loops

```
while cond do
  compute();
```

For loops

```
for indices in iteratable-expr do
  compute();
```

```
select key {
  when value1 do compute1();
  when value2 do compute2();
  otherwise do compute3();
}
```

CONTROL FLOW: BRACES VS. KEYWORDS

Of course, since compound statements are single statements, the two forms can be mixed...

Conditional statements

```
if cond then { computeA(); } else { computeB(); }
```

While loops

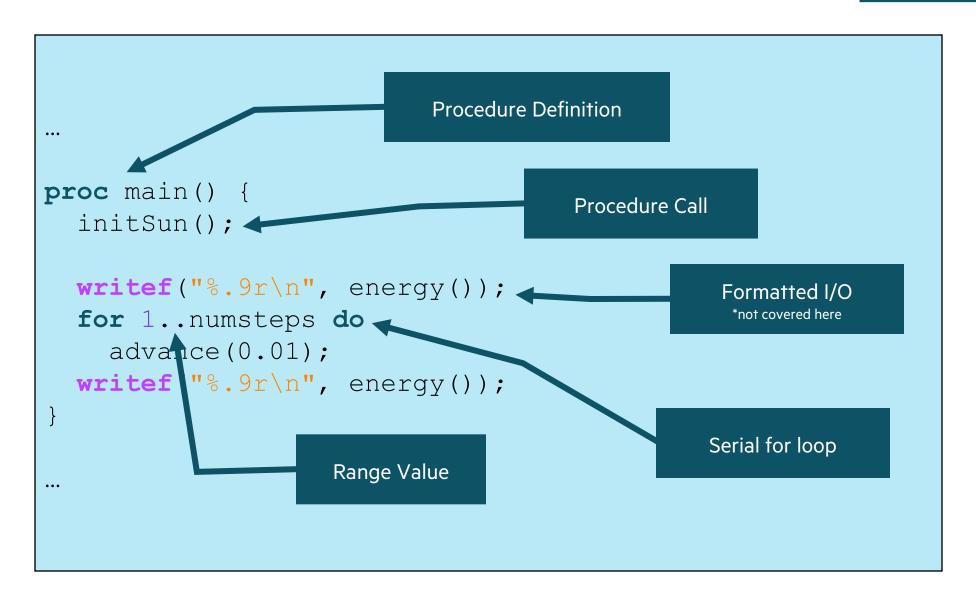
```
while cond do {
  compute();
}
```

For loops

```
for indices in iteratable-expr do {
  compute();
}
```

```
select key {
  when value1 do { compute1(); }
  when value2 do { compute2(); }
  otherwise do { compute3(); }
}
```

PROCEDURES AND ITERATORS



```
advance (0.01);
proc advance(dt) {
  for i in 1..numbodies {
    for j in i+1..numbodies {
      const dpos = bodies[i].pos - bodies[j].pos,
            mag = dt / sqrt(sumOfSquares(dpos))**3;
      bodies[i].v -= dpos * bodies[j].mass * mag;
      bodies[j].v += dpos * bodies[i].mass * mag;
  for b in bodies do
   b.pos += dt * b.v;
```

```
m_1 \mathbf{a}_1 = \frac{Gm_1m_2}{r_{12}^3} (\mathbf{r}_2 - \mathbf{r}_1) Sun-Earth
advance (0.01);
                                    m_2 \mathbf{a}_2 = \frac{Gm_1m_2}{r_{21}^3} (\mathbf{r}_1 - \mathbf{r}_2) Earth-Sun
proc advance(dt) {
  for i in 1..numbodies {
     for j in i+1..numbodies {
        const dpos = bodies[i].pos - bodies[j].pos,
                 mag = dt / sqrt(sumOfSquares(dpos)) **3;
       bodies[i].v -= dpos * bodies[j].mass * mag;
       bodies[j].v += dpos * bodies[i].mass * mag;
  for b in bodies do
     b.pos += dt * b.v;
```

5-BODY IN CHAPEL: ADVANCE()

```
advance (0.01);
                                      Procedure call
proc advance(dt) { <--</pre>
  for i in 1..numbodies {
                                        Procedure definition
    for j in i+1..numbodies {
      const dpos = bodies[i].pos - bodies[j].pos,
             mag = dt / sqrt(sumOfSquares(dpos))**3;
      bodies[i].v -= dpos * bodies[j].mass * mag;
      bodies[j].v += dpos * bodies[i].mass * mag;
  for b in bodies do
    b.pos += dt * b.v;
```

• Example to compute the area of a circle

```
proc area(radius: real): real {
  return 3.14 * radius**2;
}
writeln(area(2.0)); // 12.56
```

```
proc area(radius) {
  return 3.14 * radius**2;
}

Argument and return
```

types can be omitted

• Example of argument default values, naming

ARGUMENT INTENTS

Arguments can optionally be given intents

- (blank): varies with type; follows principle of least surprise
 - -most types: const in or const ref
 - -sync/single vars, atomics: ref
- **ref**: formal is a reference back to the actual
- const [ref | in]: disallows modification of the formal
- param/type: actual must be a param/type
- in: initializes formal using actual; permits formal to be modified
- out: copies formal into actual at procedure return
- inout: does both of the above

ARGUMENT INTENTS, BY EXAMPLE

• For some types, argument intents are needed so as to avoid inadvertent races

- Arguments can optionally be given intents.
- 'ref' intent means the actual being passed in will be modified

• Can't pass a 'const' to a 'ref' intent

- Can pass a 'const' to a 'const ref' intent
- However, can't write to a formal coming in as 'const' intent

Can't pass 'const' and 'var' into 'param' intents

```
proc foo(param x: real, type t) {
    ...
    ...
}

const r: real,
    A: [1..3] real;

// foo(r, A); // illegal: can't pass vars and consts to params and types
writeln((r, A)); // writes(0.0, [0.0, 0.0, 0.0])
```

• Can pass a literal, param, or a type into 'param' intent

```
proc foo(param x: real, type t) {
    ...
    ...
}

const r: real,
    A: [1..3] real;

foo(1.2, r.type); // OK: passing a literal/param and a type
writeln((r, A)); // writes(0.0, [0.0, 0.0, 0.0])
```

• 'in' intents cause the actual argument value to be copied into the formal

ARGUMENT INTENTS, BY EXAMPLE

• 'out' intents cause the formal value to be copied into actual argument upon return from procedure

'inout' intent is a combination of 'in' and 'out' intent

```
proc advance(dt) {
  for i in 1..numbodies {
    for j in i+1..numbodies {
      const dpos = bodies[i].pos - bodies[j].pos,
             mag = dt / sqrt(sumOfSquares(dpos))**3;
      bodies[i].v -= dpos * bodies[j].mass * mag;
      bodies[j].v += dpos * bodies[i].mass * mag;
  for b in bodies do
   b.pos += dt * b.v;
```

5-BODY IN CHAPEL: ALTERNATIVE USING ITERATORS



nbody.chpl

Use of iterator proc advance(dt) { for (i,j) in triangle(numbodies) { const dpos = bodies[i].pos - bodies[j].pos, mag = dt / sqrt(sumOfSquares(dpos)) **3; iter triangle(n) { Definition of iterator for i in 1..n do for j in i+1..n do yield (i, j);

5-BODY IN CHAPEL: ADVANCE() USING ITERATORS

```
proc advance(dt) {
  for (i, j) in triangle(numbodies) {
    const dpos = bodies[i].pos - bodies[j].pos,
          mag = dt / sqrt(sumOfSquares(dpos)) **3;
    bodies[i].v -= dpos * bodies[j].mass * mag;
    bodies[j].v += dpos * bodies[i].mass * mag;
  for b in bodies do
   b.pos += dt * b.v;
```

HANDS ON: WHERE MIGHT WE CONSIDER PARALLELIZING N-BODY

Look at 'nbody.chpl' and identify...



- 'for' loops that can be parallelized
- 'for' loops that need to stay serial to keep meaning
- 'for' loops that are "mostly" parallel but have something like +=

Can be parallelized

Inherently serial loop

Can be parallelized but have to avoid races when adding into velocity field

```
for b in bodies do
 b.pos += dt * b.v;
for 1...numsteps do
  advance (0.01);
for i in 1..numbodies {
  for j in i+1..numbodies {
    const dpos = bodies[i].pos - bodies[j].pos,
          mag = dt / sqrt(sumOfSquares(dpos))**3;
    bodies[i].v -= dpos * bodies[j].mass * mag;
    bodies[j].v += dpos * bodies[i].mass * mag;
```

OUTLINE: CHAPEL BASICS

- Running Example: n-body computation (Hands On)
- Variables, Constants, and Operators
- Records and Classes
- Tuples
- Arrays
- Writing out Tuples, Records, and Arrays (Hands On)
- Main() Procedure
- Ranges and basic control flow
- Procedures and iterators
- Where might we parallelize the n-body computation? (Hands On)

LEARNING OBJECTIVES FOR TODAY'S CHAPEL TUTORIAL

- Familiarity with the Chapel execution model including how to run codes in parallel on a single node, across nodes, and both
- Learn Chapel concepts by compiling and running provided code examples
 - ✓ Serial code using map/dictionary, (k-mer counting from bioinformatics)
 - ✓ Parallelism and locality in Chapel
 - ✓ Distributed parallelism and 1D arrays, (processing files in parallel)
 - ✓ Chapel basics in the context of an n-body code
 - Distributed parallelism and 2D arrays, (heat diffusion problem)
 - How to parallelize histogram
 - Using CommDiagnostics for counting remote reads and writes
 - Chapel and Arkouda best practices including avoiding races and performance gotchas
- Where to get help and how you can participate in the Chapel community



ONE DAY CHAPEL TUTORIAL

- 9-10:30: Getting started using Chapel for parallel programming
- 10:30-10:45: break
- 10:45-12:15: Chapel basics in the context of the n-body example code
- 12:15-1:15: lunch
- 1:15-2:45: Distributed and shared-memory parallelism especially w/arrays (data parallelism)
- 2:45-3:00: break
- 3:00-4:30: More parallelism including for asynchronous parallelism (task parallelism)
- 4:30-5:00: Wrap-up including gathering further questions from attendees

CHAPEL RESOURCES

Chapel homepage: https://chapel-lang.org

• (points to all other resources)

Social Media:

• Twitter: <a>@ChapelLanguage

Facebook: @ChapelLanguage

• YouTube: http://www.voutube.com/c/ChapelParallelProgrammingLanguage

Community Discussion / Support:

Discourse: https://chapel.discourse.group/

Gitter: https://gitter.im/chapel-lang/chapel

Stack Overflow: https://stackoverflow.com/questions/tagged/chapel

• GitHub Issues: https://github.com/chapel-lang/chapel/issues



What is Chapel? What's New? Upcoming Events
Job Opportunities

How Can I Learn Chapel? Contributing to Chapel

Performance Powered by Chapel

User Resources Developer Resources

Social Media / Blog Posts

Contributors / Credits chapel_info@cray.com







What is Chapel?

Chapel is a programming language designed for productive parallel computing at scale.

The Chapel Parallel Programming Language

Why Chapel? Because it simplifies parallel programming through elegant support for:

- · distributed arrays that can leverage thousands of nodes' memories and cores
- a global namespace supporting direct access to local or remote variables
- . data parallelism to trivially use the cores of a laptop, cluster, or supercomputer
- · task parallelism to create concurrency within a node or across the system

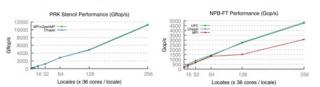
Chapel Characteristics

- productive: code tends to be similarly readable/writable as Python
- · scalable: runs on laptops, clusters, the cloud, and HPC systems
- fast: performance competes with or beats C/C++ & MPI & OpenMP
- · portable: compiles and runs in virtually any *nix environment
- · open-source: hosted on GitHub, permissively licensed

New to Chapel?

As an introduction to Chapel, you may want to...

- · watch an overview talk or browse its slides
- read a blog-length or chapter-length introduction to Chapel
- · learn about projects powered by Chapel
- check out performance highlights like these:



· browse sample programs or learn how to write distributed programs like this one:

```
// use the Cyclic distribution library
                         // use --n=<val> when executing to override this default
forall i in {1..n} dmapped Cyclic(startIdx=1) do
 writeln("Hello from iteration ", i, " of ", n, " running on node ", here.id);
```