



# CHAPEL BASICS, PART II

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Chapel Team, edited by Michelle Strout

April 8, 2025

# HANDS ON: HOW TO DO THE HANDS ON



01-hello.chpl

## Example codes for Chapel tutorial slides

- <https://github.com/UofA-CSc-372-Spring-2025/CSc372Spring2025-CourseMaterials/tree/main/Sandboxes/ChapelTutorialExamples>

## Using a container on your laptop

- First, install docker for your machine and start it up (see the README.md for more info)
- Then, use the chapel-gasnet docker container

```
docker pull docker.io/chapel/chapel-gasnet      # takes about 5 minutes
cd CSc372Spring2025-CourseMaterials/Sandboxes/ChapelTutorialExamples/
docker run --rm -it -v "$PWD":/workspace chapel/chapel-gasnet
root@589405d07f6a:/opt/chapel# cd /workspace
root@xxxxxxxx:/myapp# chpl 01-hello.chpl
root@xxxxxxxx:/myapp# ./01-hello -nl 1
```



# PLAN

- **Announcements**

- Everyone should have their Final Project assignments (go over final project expectations)
- SA7 is due Friday April 11th

- **Last time**

- TopHat Questions
- ICA10 Quiz
- Chapel Programming Basics in the context of an Nbody simulation, part I

- **Today**

- Chapel programming basics in the context of an Nbody simulation, part II

# SA7 EDITS FROM YESTERDAY

- See piazza post, screen shot below for reference

Actions ▾

## SA7 Chapel edits

The SA7 assignment writeup, <https://github.com/UofA-CSc-372-Spring-2025/CSc372Spring2025-CourseMaterials/blob/main/SmallAssignmentWriteups/sa7-chapel.md>, and initial files, <https://github.com/UofA-CSc-372-Spring-2025/sa7-chapel-start>, have undergone some modifications.

- The array of files is being tested in sa7-student-tests.chpl using a set instead of just comparing the array values since sometimes the order didn't match.
- The rgb to gray conversion and the sobel edge detector descriptions have been clarified in the writeup. They gray value needs to be put in all of the rgb channels, before and after edge detection.
- There is are some tests for the values in the histogram now.
- The Gradescope autograder has been set up as of Monday April 7th at 8:15pm. Already submitted assignments have been regraded. The grading tests are somewhat different that the tests provided in sa7-student-tests.chpl.

## OUTLINE: OVERVIEW OF PROGRAMMING IN CHAPEL

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- Main( ) Procedure
- Ranges and basic control flow
- Procedures and iterators
- Where might we parallelize the n-body computation? (Hands On)



MAIN() PROCEDURE

## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

...

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

...

## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

...

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

...

Procedure Definition



## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

...

```
proc main() {  
  initSun();
```

Procedure Call

```
  writef("%.9r\n", energy());
```

```
  for 1..numsteps do
```

```
    advance(0.01);
```

```
  writef("%.9r\n", energy());
```

```
}
```

...

## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

Activity: Using the table at <https://chapel-lang.org/docs/modules/standard/IO/FormattedIO.html>, format the energy values in three different ways.

...

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

...

Formatted I/O

## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

```
...  
  
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  
     $\text{low} \leq \text{high}$ :  $\text{low}, \text{low}+1, \text{low}+2, \dots, \text{high}$   
     $\text{low} > \text{high}$ : degenerate (an empty range)  
     $\text{low}$  or  $\text{high}$  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

 02-range-operators.chpl

```
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
```

**Activity:** Experiment with 02-range-operators.chpl. What are the two different ways the pound sign (#) are being used?

## 5-BODY IN CHAPEL: MAIN( )



nbody.chpl

```
...  
  
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]); }           // DO RE MI  
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 statements

```
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 iterable-expr {  
    compute();  
}
```

- Select statements

```
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 iterable-expr do  
    compute();
```

- Select statements

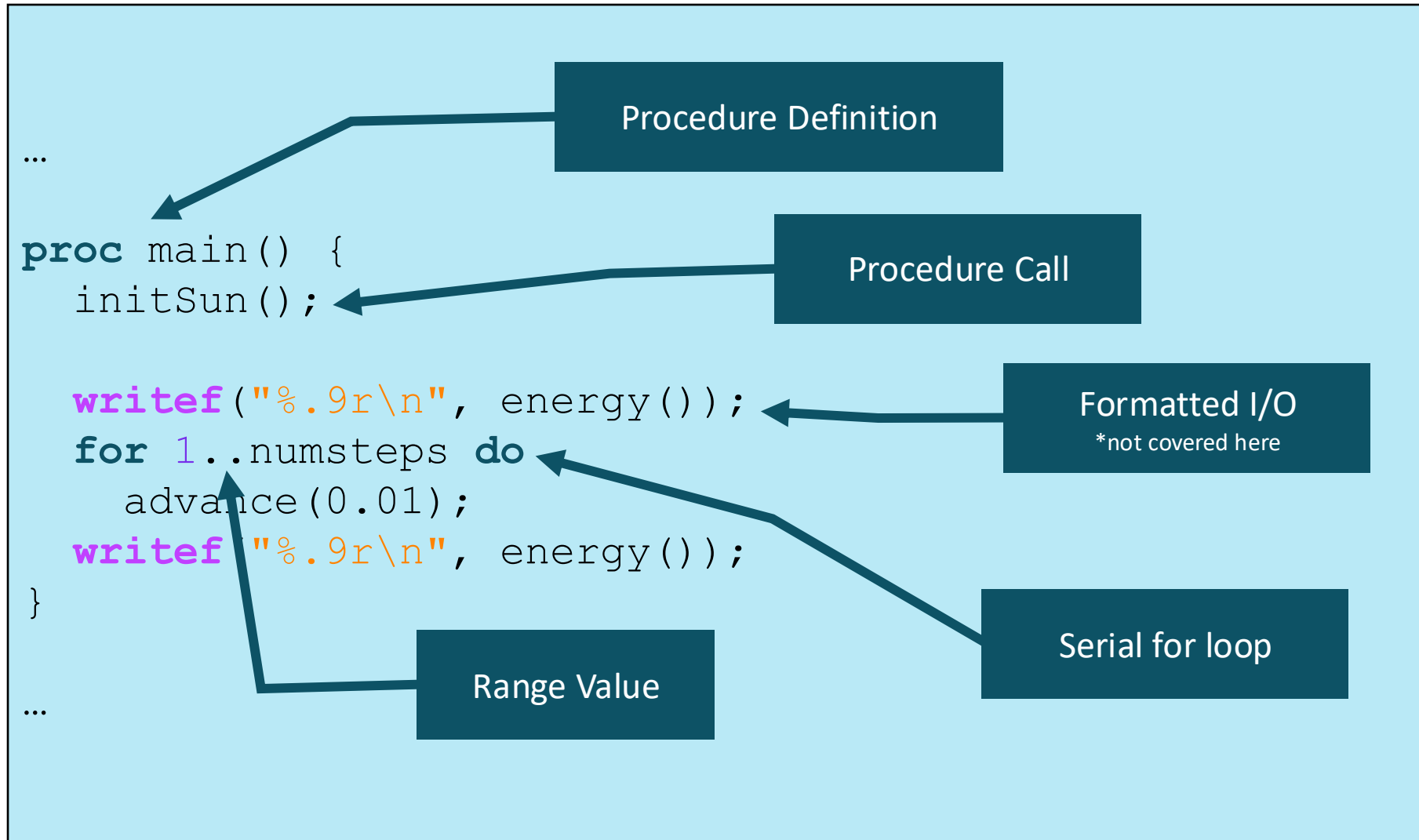
```
select key {  
    when value1 do compute1();  
    when value2 do compute2();  
    otherwise   do compute3();  
}
```

# PROCEDURES AND ITERATORS

# 5-BODY IN CHAPEL: MAIN( )



nbody.chpl



## 5-BODY IN CHAPEL: ADVANCE( )



nbody.chpl

```
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;  
}
```

## 5-BODY IN CHAPEL: ADVANCE()



nbody.chpl

```
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;
```

```
}
```

$$m_1 \mathbf{a}_1 = \frac{G m_1 m_2}{r_{12}^3} (\mathbf{r}_2 - \mathbf{r}_1) \quad \text{Sun-Earth}$$

$$m_2 \mathbf{a}_2 = \frac{G m_1 m_2}{r_{21}^3} (\mathbf{r}_1 - \mathbf{r}_2) \quad \text{Earth-Sun}$$



## 5-BODY IN CHAPEL: ADVANCE()



nbody.chpl

```
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;  
}
```

Procedure call

Procedure definition

# PROCEDURES, BY EXAMPLE

- 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

```
proc writeCoord(x: real = 0.0, y: real = 0.0) {  
    writeln((x, y));  
}
```

```
writeCoord(2.0);           // (2.0, 0.0)  
writeCoord(y=2.0);         // (0.0, 2.0)  
writeCoord(y=2.0, 3.0);    // (3.0, 2.0)
```

# 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

**Question: Why might it be useful for the compiler to have this information?**



# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(x: real, y: [] real) {  
    // x = 1.2;    // illegal: scalars are passed 'const in' by default  
    // y = 3.4;    // illegal: 'ref' by default for arrays is deprecated  
}  
  
var r: real,  
    A: [1..3] real;  
  
foo(r, A);  
  
writeln((r, A));
```

# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(ref x: real, ref y: [] real) {  
  x = 1.2;  // OK: actual is modified  
  y = 3.4;  // OK: actual is modified  
}  
  
var r: real,  
    A: [1..3] real;  
  
foo(r, A);  
  
writeln((r, A));  // writes (1.2, [3.4, 3.4, 3.4])
```

# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(ref x: real, ref y: [] real) {  
  x = 1.2; // OK: actual is modified  
  y = 3.4; // OK: actual is modified  
}  
  
const r: real,  
      A: [1..3] real;  
  
// foo(r, A); // illegal, can't pass a constant to a 'ref' intent  
  
writeln((r, A)); // writes (0.0, [0.0, 0.0, 0.0])
```

# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(const ref x: real, const ref y: [] real) {  
    // x = 1.2;    // illegal: can't modify constant arguments  
    // y = 3.4;    // illegal: can't modify constant arguments  
}  
  
const r: real,  
      A: [1..3] real;  
  
foo(r, A);    // OK to create constant references to constants  
  
writeln((r, A));    // writes (0.0, [0.0, 0.0, 0.0])
```

**Question: Why would we want to pass something by reference if we didn't plan on modifying it?**

# ARGUMENT INTENTS, BY EXAMPLE

- 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])
```



# ARGUMENT INTENTS, BY EXAMPLE

- 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])
```

# ARGUMENT INTENTS, BY EXAMPLE

- 'in' intents cause the formal variable to get its own value of the actual argument

```
proc foo(in x: real, in y: [] real) {  
  x = 1.2;  // OK: local copy is modified  
  y = 3.4;  // OK: local copy is modified  
}  
  
var r: real,  
    A: [1..3] real;  
  
foo(r, A);  
  
writeln((r, A));  // writes (0.0, [0.0, 0.0, 0.0])
```

# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(out x: real, out y: [] real) {  
  x = 1.2;  // OK: local copy is modified  
  y = [3.4, 3.4, 3.4];  // OK: local copy is modified  
}  
  
var r: real,  
    A: [1..3] real;  
  
foo(r, A);  
  
writeln((r, A));  // writes (1.2, [3.4, 3.4, 3.4])
```

# ARGUMENT INTENTS, BY EXAMPLE

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

```
proc foo(inout x: real, inout y: [] real) {  
    x = 1.2; // OK: local copy is modified  
    y = 3.4; // OK: local copy is modified  
}  
  
var r: real,  
    A: [1..3] real;  
  
foo(r, A);  
  
writeln((r, A)); // writes (1.2, [3.4, 3.4, 3.4])
```

## 5-BODY IN CHAPEL: ADVANCE( )



nbody.chpl

```
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;  
    ...  
  }  
  ...  
}
```

Definition of iterator

```
iter triangle(n) {  
  for i in 1..n do  
    for j in i+1..n do  
      yield (i,j);  
    }  
}
```

## 5-BODY IN CHAPEL: ADVANCE( ) USING ITERATORS



nbody.chpl

```
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 +=

See <https://chapel-lang.org/docs/technotes/reduceIntents.html>

Can be parallelized

```
for b in bodies do  
  b.pos += dt * b.v;
```

Inherently serial loop

```
for 1..numsteps do  
  advance(0.01);
```

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

```
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;  
  }  
}
```



# CHAPEL RESOURCES

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

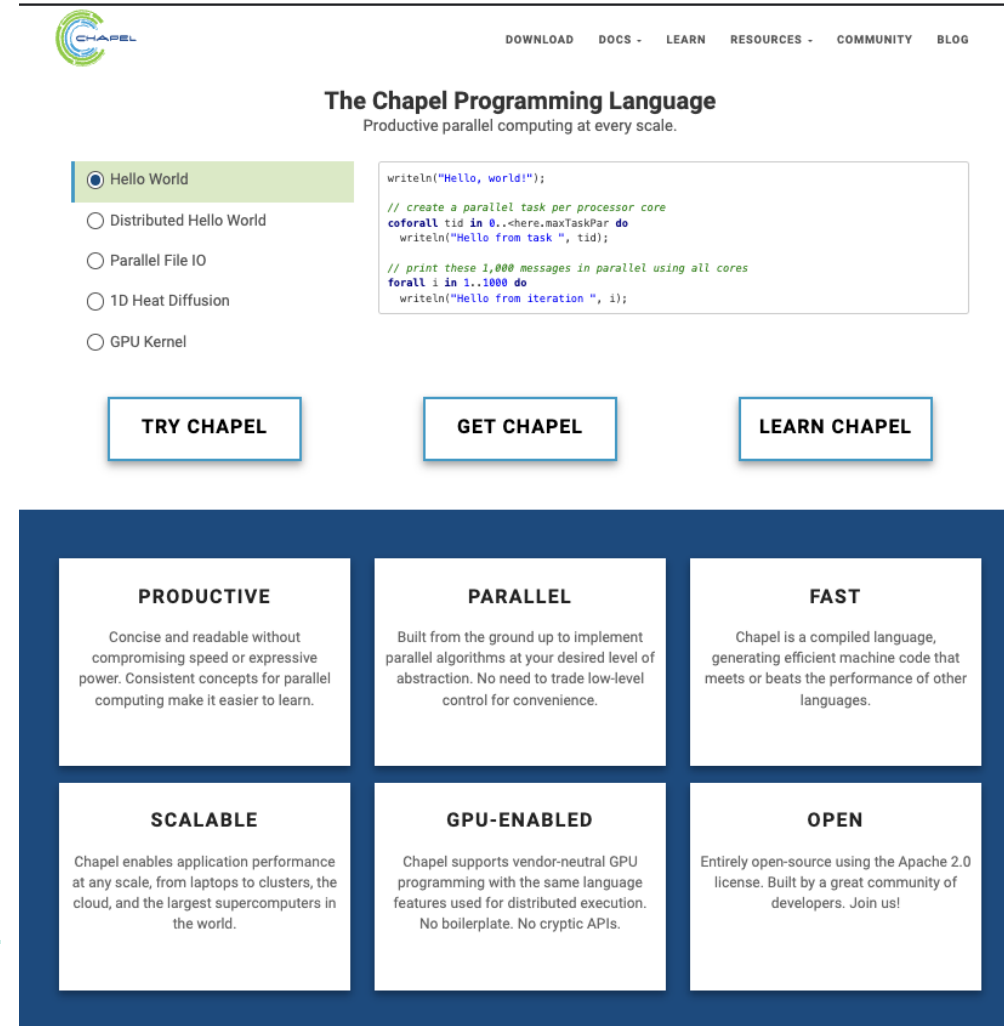
- (points to all other resources)

## Social Media:

- Twitter: [@ChapelLanguage](https://twitter.com/ChapelLanguage)
- Facebook: [@ChapelLanguage](https://facebook.com/ChapelLanguage)
- YouTube: <http://www.youtube.com/c/ChapelParallelProgrammingLanguage>

## Community Discussion / Support:

- Discord: <https://discord.com/invite/xu2xg45yqH>
- Stack Overflow: <https://stackoverflow.com/questions/tagged/chapel>
- GitHub Issues: <https://github.com/chapel-lang/chapel/issues>



The screenshot shows the Chapel Programming Language homepage. At the top, there's a navigation bar with links: DOWNLOAD, DOCS, LEARN, RESOURCES, COMMUNITY, and BLOG. The main heading is "The Chapel Programming Language" with the tagline "Productive parallel computing at every scale." Below this, there's a section for "Hello World" with a radio button selected, and four other options: Distributed Hello World, Parallel File IO, 1D Heat Diffusion, and GPU Kernel. To the right of these options is a code editor showing a "Hello, world!" program. Below the code editor are three buttons: TRY CHAPEL, GET CHAPEL, and LEARN CHAPEL. At the bottom, there's a grid of six boxes, each describing a feature of Chapel: PRODUCTIVE, PARALLEL, FAST, SCALABLE, GPU-ENABLED, and OPEN.

**PRODUCTIVE**  
Concise and readable without compromising speed or expressive power. Consistent concepts for parallel computing make it easier to learn.

**PARALLEL**  
Built from the ground up to implement parallel algorithms at your desired level of abstraction. No need to trade low-level control for convenience.

**FAST**  
Chapel is a compiled language, generating efficient machine code that meets or beats the performance of other languages.

**SCALABLE**  
Chapel enables application performance at any scale, from laptops to clusters, the cloud, and the largest supercomputers in the world.

**GPU-ENABLED**  
Chapel supports vendor-neutral GPU programming with the same language features used for distributed execution. No boilerplate. No cryptic APIs.

**OPEN**  
Entirely open-source using the Apache 2.0 license. Built by a great community of developers. Join us!