



# Developing Stencil Code using the YASK Framework

This tutorial applies to YASK version 4.00 (and later; see README.md for exceptions)

Chuck Yount Intel Americas, Inc.

Nov. 8, 2022 © 2019-2022 Intel Corporation

# **NOTICES AND DISCLAIMERS**

Intel technologies' features and benefits depend on system configuration and may require enabled hardware, software or service activation. Performance varies depending on system configuration.

No computer system can be absolutely secure.

Tests document performance of components on a particular test, in specific systems. Differences in hardware, software, or configuration will affect actual performance. For more complete information about performance and benchmark results, visit <a href="http://www.intel.com/benchmarks">http://www.intel.com/benchmarks</a>.

Software and workloads used in performance tests may have been optimized for performance only on Intel microprocessors. Performance tests, such as SYSmark and MobileMark, are measured using specific computer systems, components, software, operations and functions. Any change to any of those factors may cause the results to vary. You should consult other information and performance tests to assist you in fully evaluating your contemplated purchases, including the performance of that product when combined with other products. For more complete information visit <a href="http://www.intel.com/benchmarks">http://www.intel.com/benchmarks</a>.

Intel® Advanced Vector Extensions (Intel® AVX)\* provides higher throughput to certain processor operations. Due to varying processor power characteristics, utilizing AVX instructions may cause a) some parts to operate at less than the rated frequency and b) some parts with Intel® Turbo Boost Technology 2.0 to not achieve any or maximum turbo frequencies. Performance varies depending on hardware, software, and system configuration and you can learn more at <a href="http://www.intel.com/go/turbo">http://www.intel.com/go/turbo</a>.

Intel's compilers may or may not optimize to the same degree for non-Intel microprocessors for optimizations that are not unique to Intel microprocessors. These optimizations include SSE2, SSE3, and SSSE3 instruction sets and other optimizations. Intel does not guarantee the availability, functionality, or effectiveness of any optimization on microprocessors not manufactured by Intel. Microprocessor-dependent optimizations in this product are intended for use with Intel microprocessors. Certain optimizations not specific to Intel microarchitecture are reserved for Intel microprocessors. Please refer to the applicable product User and Reference Guides for more information regarding the specific instruction sets covered by this notice.

Cost reduction scenarios described are intended as examples of how a given Intel-based product, in the specified circumstances and configurations, may affect future costs and provide cost savings. Circumstances will vary. Intel does not guarantee any costs or cost reduction.

Intel does not control or audit third-party benchmark data or the web sites referenced in this document. You should visit the referenced web site and confirm whether referenced data are accurate. Intel, the Intel logo, and Intel Xeon are trademarks of Intel Corporation in the U.S. and/or other countries.

\*Other names and brands may be claimed as property of others.

© 2018 Intel Corporation.



### **Outline**

#### Introduction

- What is YASK?
- Motivation and example YASK application
- Scope of this presentation

#### Basic features and usage

- Download, build, and test
- High-level flow
- Writing a simple stencil in the DSL
- Building the YASK library
- Testing and tuning your stencil
- Multi-process usage via MPI

### Using the APIs

- Access, terms, and basic usage
- Advanced APIs and exceptions

#### Advanced stencils and tuning

- Multiple stencils
- Boundary regions (sub-domains)
- Temporal conditions
- Vector folding
- Cluster, bundles, and stages
- FP precision and prefetching
- 3 layers of temporal tiling
- Nested OpenMP in sub-blocks
- More on the auto-tuners

#### **GPU Offloading**

#### Wrap-up

- Current work
- Further reading and call to action

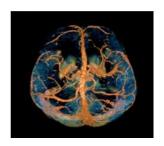


# INTRODUCTION

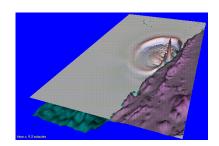
### YASK: "Yet Another Stencil Kit"

YASK is a software *framework* for the rapid development of HPC stencil-based applications

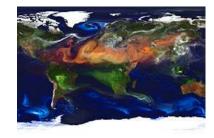
- Stencil: an iterative kernel that updates elements in one or more N-dimensional vars using a fixed pattern of computation on neighboring elements
- Fundamental algorithm in many scientific simulations, e.g., finite-difference-method (FDM) approximations of differential equations describing various physical phenomena



**Image Processing** 



Seismic Modeling



Weather Simulation

Images from <a href="https://commons.wikimedia.org">https://commons.wikimedia.org</a>

#### About the name:

- "YASK" is named in the tradition of Linux utilities such as YACC, YAML and YAKL: <a href="https://en.wikipedia.org/wiki/Yet\_another">https://en.wikipedia.org/wiki/Yet\_another</a>
- Prior to version 3, "YASK" stood for "Yet Another Stencil Kernel"

### Technical and business motivation

#### Rapid Development

- Stencils in YASK are coded in a simple DSL (domain-specific language)
  - YASK programmer only needs to describe what to do, not how to do it
  - YASK compiler generates high-perf code from the DSL description
  - Can easily and quickly change the stencil and generate new code
  - Supports arbitrary dimensions, complex stencils, boundary conditions, and more
- Can easily and quickly try different tuning features and parameters without recoding
  - Many complex optimization techniques are available immediately
  - Supports cluster scaling, spatial and temporal tiling, vector folding, and more
- Generated code compiles into a library with documented C++ and Python\* APIs to facilitate integration into real HPC applications

#### Performance Portability

- Can re-target stencils for different Intel® CPU and GPU platforms by generating multiple libraries from single DSL description
- Future YASK features and supported platforms can be leveraged immediately without recoding



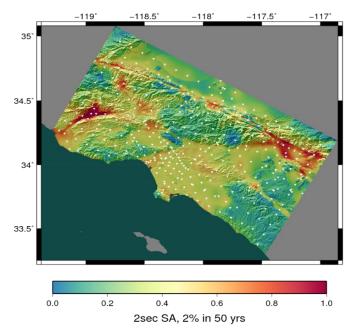
### Example application of YASK

### AWP-ODC: Anelastic Wave Propagation-Olsen, Day, Cui

- Software that simulates seismic wave propagation after a fault rupture
- Widely used by the Southern California Earthquake Center (SCEC) community

#### **AWP-ODC-OS**

- First ever open-source release in 2016 (BSD-2 license), including port to Intel Xeon Phi processor, under development by San Diego Supercomputer Center (SDSC) at Univ. of CA, San Diego (UCSD)
- Demonstrated on >9000 nodes of Cori supercomputer



- CyberShake Study 15.4 hazard map for 336 sites around Southern California
- · Warm colors represent areas of high hazard

Content on this slide courtesy of UCSD



### Scope of this presentation

#### Goals

- Give a practical introduction to using YASK—a tutorial
- Provide an overview of the major YASK performance techniques
- Deliver enough information for someone already familiar with the application of stencil codes to start using YASK

#### Non-goals

- Not providing a tutorial on
  - Finite-difference methods, seismic modeling, brain imaging, etc.
  - C++, Linux\*, OpenMP\*, MPI\*, github\*, etc.
  - Intel® instruction sets, CPU/GPU architecture, etc.
- Not explaining in-depth how the YASK performance techniques work
- Not providing performance data on all the various trade-offs
- Not describing the internal software architecture
- Not a reference manual for the APIs



# BASIC FEATURES AND USAGE

### Download, build, and test

#### Code access

- Download from Intel's github\* project
  - git clone https://github.com/intel/yask
  - MIT open-source license
- Builds and runs are made from the top-level directory: cd yask

#### **Pre-requisites**

- See README.md for complete list
- Only supported OS is Linux (no specific distribution recommended)
- Intel® C++ compiler needed for performance
  - Install "Intel® oneAPI HPC Toolkit"
  - https://www.intel.com/content/www/us/en/developer/tools/oneapi/toolkits.html
- Common Linux utilities (gmake, perl, awk, python, etc.)





### Example 1: Iso3dfd stencil

### Description

- Isotropic 3D finite-difference code found in seismicimaging software used by energy-exploration companies to predict the location of oil and gas deposits
  - Simple stencil with only one updated variable

### Recipes for building and running on CPU

- Makefile and run script will automatically determine architecture
  - make clean; make -j stencil=iso3dfd
  - bin/yask.sh -stencil iso3dfd -g 1024
- The -g 1024 option sets the global-domain size to a cube, 1024 elements on a side



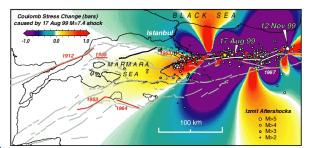




### Example 2: AWP stencil

#### Description

- Primary compute kernel for earthquake simulator described in the introduction
  - More complex problem that consists of 26 vars in a staggeredgrid formulation



### Recipes for building and running on CPU

Makefile and run script will automatically determine architecture



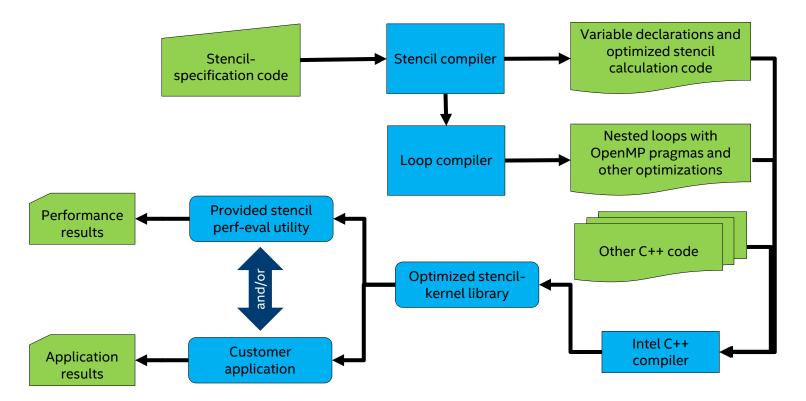
- bin/yask.sh -stencil awp -gx 1024 -gy 1024 -gz 128
- The -g\* options set the global-domain size to specific sizes in each dimension
  - Compare this to the simpler -g option in the iso3dfd example that set the global-domain size in all three spatial dimensions
  - This is a common technique used for all spatial-dimension command-line options

Image from https://commons.wikimedia.org/wiki/File:Izmit\_11-12-99.gif. Public domain--U.S.G.S.

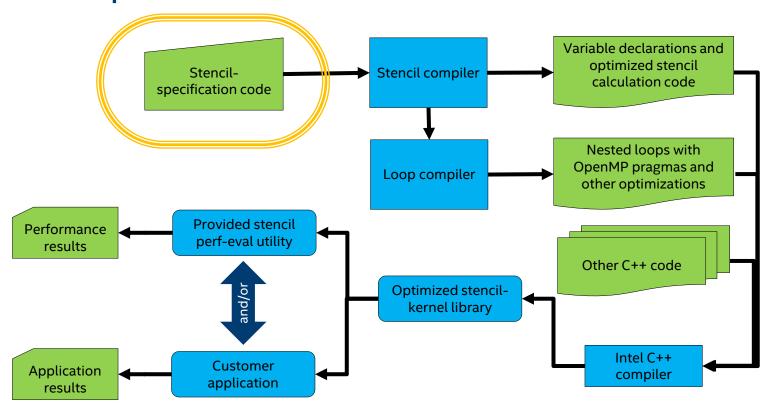




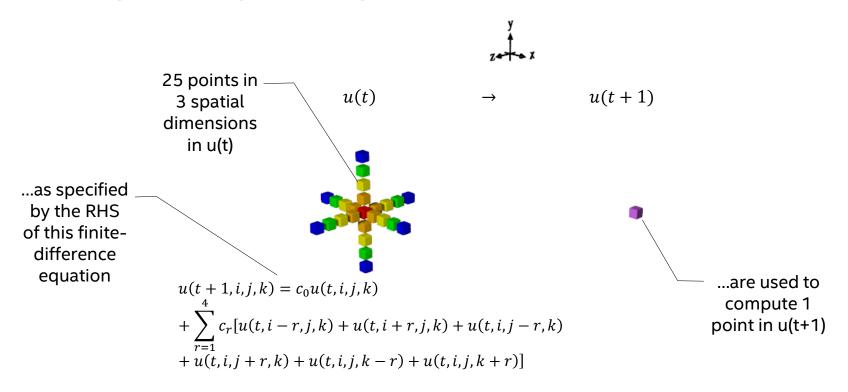
# High-level tool-chain flow



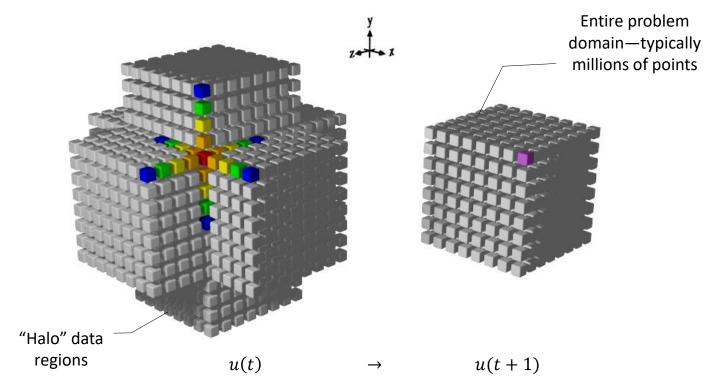
# Stencil specification



# Example simple 25-point 3-D stencil



# Stencil will be applied over entire problem domain



Repeat for u(t+2), u(t+3), ...



# Example stencil DSL code

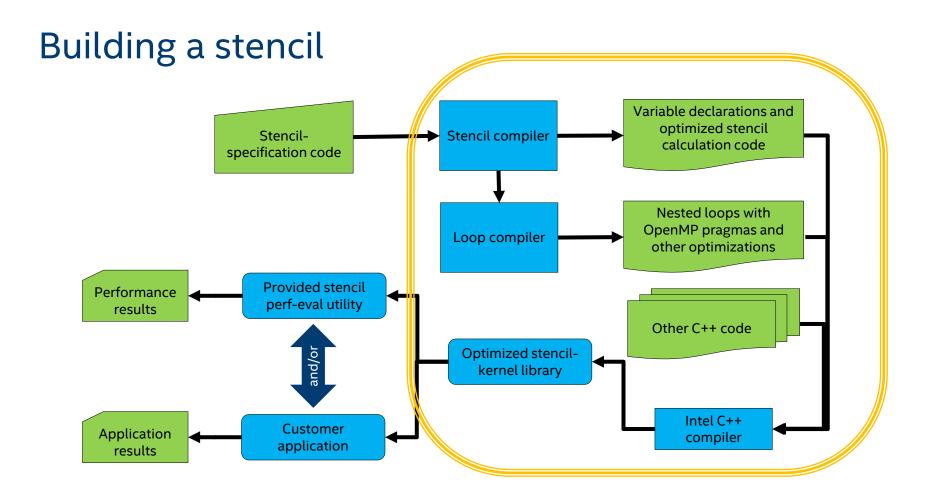
Mathematical description of finite-difference approximation implemented in DSL code



```
u(t + 1, i, j, k) = c_0 u(t, i, j, k)
+ \sum_{r=1}^{4} c_r [u(t, i - r, j, k) + u(t, i + r, j, k) + u(t, i, j - r, k) + u(t, i, j + r, k) + u(t, i, j, k - r) + u(t, i, j, k + r)]
```

```
yc solution base (or
                                                       yc solution with radius base to
#include "yask compiler api.hpp"
using namespace yask;
                                                               get radius parameter)
namespace {
class MyStencil: public yc solution with radius base
                                                            Overload define() method
public:
 MyStencil(int radius=4) :
                                                             Declare 1 time and 3 space
      yc solution with radius base ("my stencil", radius)
                                                               indices and 4D "u" var
 virtual void define()
    auto t = new step index("t");
    auto x = new domain index("x");
                                                              Declare misc index and 1D
    auto y = new domain index("y");
                                                                "c" array for coefficients
    auto z = new domain index("z");
    yc var proxy u("U", get soln(),
    auto i = new misc index("i");
                                                              Write expression for value
    yc var proxy c("C", get soln(), { i });
                                                                     at step t+1
    auto nu = c(0) * u(t, x, y, z);
    for (int r = 1; r \le get radius(); r++)
                                                           Use EQUALS operator to define
       nu += c(r) * (u(t, x-r, y, z) + u(t, x+r, y, z) +
                     u(t, x, y-r, z) + u(t, x, y+r, z)
                                                                u(t+1)—this is not an
                     u(t, x, y, z-r) + u(t, x)
                                                                    assignment!
    u(t+1, x, y, z) EQUALS nu;
};
                                                          Register this stencil with the YASK
                                                                   compiler utility
static MyStencil MyStencil instance;
```

Derive new class from



# Building your example stencil

#### Build your stencil and the YASK compiler

- Put the code from a couple of slides back into a new file src/stencils/MyStencil.cpp
  - If you really don't want to copy the example stencil during this tutorial, just substitute iso3dfd for my\_stencil in the
    example commands from here on
  - If you want to write your own stencil, see the examples in src/stencils and the full DSL documentation for the YASK Compiler as discussed in the "Using the APIs" section later in this tutorial
- Optional: rebuild and run the YASK compiler manually
  - Build the compiler: make -j compiler
  - Run the compiler: bin/yask compiler.exe -h
    - "my stencil" should be listed as a valid parameter to the -stencil option
    - Lots of other options shown; most of these are passed automatically from the Makefile

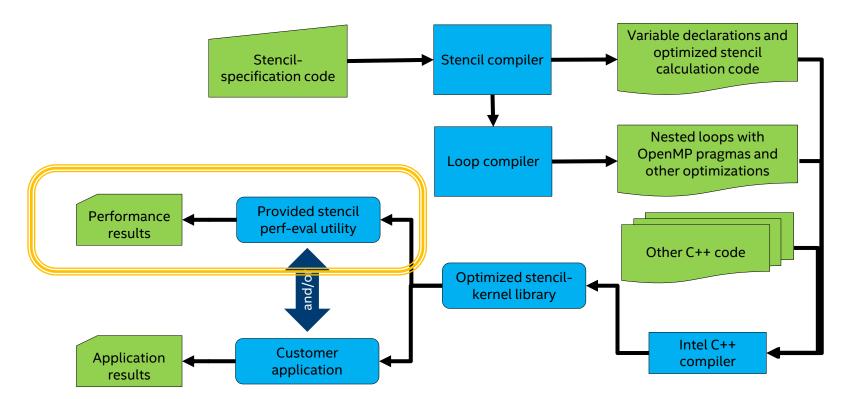
#### Build the YASK kernel for a CPU target

- Build the kernel: make -j stencil=my\_stencil radius=4
  - Targets the instruction-set architecture of your current platform by default
    - If you want to cross-compile, specify arch=isa-code, e.g., avx2
    - Targeting a GPU device will be covered later
  - Uses all the default parameters for static options such as data layout and prefetching
  - Dynamic options such as problem size are specified at run-time
- If you skipped the step of building the compiler above, it will be done here for you automatically

(intel<sup>2</sup>)

YASK tutorial

# Running a stencil



# Running a stencil on a CPU platform

### Run the provided benchmark utility

- bin/yask.sh -stencil my\_stencil -g 512 -no-pre\_auto\_tune
  - Problem size is specified as 512 in the 3 spatial dimensions (512 $^3$  points) using -g option ("g" for global-domain)
  - Auto-tuner is disabled; other parameters are set to their defaults
  - The script will use MPI to start a separate process on each NUMA node
    - The global-domain size is automatically decomposed across the processes
    - By default, uses shared-memory (shm) buffers for inter-process communication
      - Override communication type with -no-use shm option, which forces MPI calls between all processes
      - May also have to use this if your platform limits the amount of shm that can be allocated
  - The utility prints lots of stats about the stencil, memory usage, performance, etc.
  - Runs 3 trials for about 10 sec each and reports the best and median time and throughput (rate) stats
    - The most important performance statistic is usually "mid-throughput (num-points/sec)", which is printed near the end of the output
  - A log file is kept in the "logs" directory, which includes some compilation and platform information for posterity



# Automatically tuning the size parameters

#### Use the auto-tuner

- bin/yask.sh -stencil my stencil -g 512
  - As above, but with the block-size auto-tuner enabled (by default)
    - The block size determines the amount of work done by each OpenMP thread at any given time
    - The utility prints each block size that the auto-tuner tries and its measured performance, e.g.,

```
auto-tuner: searching block sizes...
auto-tuner: search-dist=16: 3.59 steps/sec (2 steps(s) in 556.7m secs) with size x=96 * y=28 * z=96 -- best so far auto-tuner: search-dist=16: 3.74 steps/sec (2 steps(s) in 533.7m secs) with size x=32 * y=4 * z=80 -- best so far
```

When the tuner is done, it prints the final sizes, e.g.,

```
auto-tuner: applying size t=0 * x=96 * y=4 * z=80
auto-tuner: done
Waiting for auto-tuner to converge on all ranks...
Auto-tuner done after 73.62 secs
Final settings:
auto-tuner: mega-block-size:
auto-tuner: block-size:
auto-tuner: block-size:
t=0 * x=256 * y=512 * z=512
t=0 * x=96 * y=4 * z=80
```

- Note that the tuned block size printed here is different than at the beginning of the run
  - Using other auto-tuner options, the other types of blocks (explained later) can be tuned
- Compare output and performance of the two runs with and without the auto-tuner





# Hand-tuning the size parameters

#### Block size

- From the log file of the last run, find the best block size found by the auto-tuner as highlighted on the previous slide
- Specify this manually to bypass the auto-tuner, e.g.,
  - bin/yask.sh -stencil my\_stencil -g 512 -no-pre\_auto\_tune \
     -bx 96 -by 4 -bz 80
  - You should get similar performance
  - Play around with block size to see its effect
  - Contribution opportunity: github ticket #159 to save and reuse the settings automatically

### General usage of size-parameter options

- Most domain-dimension sizes can be specified in two ways
  - Spatial dims set separately as we did with block size, e.g., -bx, -by, -bz
  - Leave off the dimension names to set all domain sizes, e.g., -b
- Play around with various problem sizes to see the effect on memory and performance





# Getting a list of options

The YASK test utility is composed of a script driver and a binary

- Print script options: bin/yask.sh -h
  - There is only a short list of options here
  - These are just the ones that control how the binary is launched from the script
- Print binary options: bin/yask.sh -stencil my\_stencil -help
  - Need to specify the stencil because there are different binaries for each
    - If you don't specify a stencil, the script will show the ones that have been compiled
  - The long list of options might be overwhelming; we'll point out the most important ones in this tutorial



# Scaling out to multiple nodes in a cluster

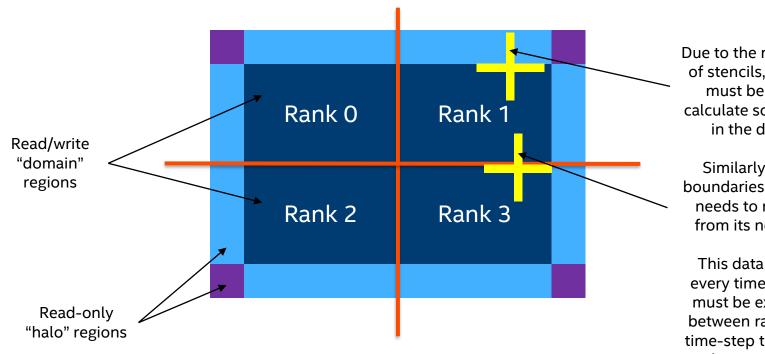
### Domain decomposition

- Many HPC problem sizes won't fit in the DRAM memory of one platform
- Most HPC problems can be divided spatially across multiple nodes in a cluster via domain decomposition

### YASK implementation

- Data exchanges use two-sided MPI\* communication
- Important: strong vs. weak HPC scaling
  - Global-domain size parameters (-g) apply to the overall problem size
    - This implements a strong-scaling model, where the overall problem size remains constant regardless of the number of MPI ranks
  - Local-domain size parameters (-1) apply to the size on each MPI rank
    - This implements a weak-scaling model, where the overall problem size increases with the number of ranks

# Conceptual view of 2D, 2×2 rank domains



Due to the read-radius of stencils, halo data must be read to calculate some points in the domain.

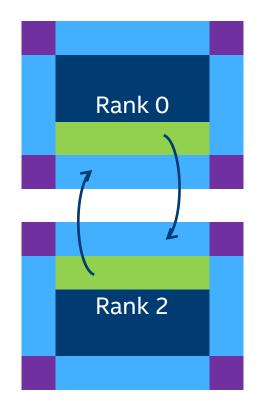
Similarly, at rank boundaries, each rank needs to read data from its neighbors.

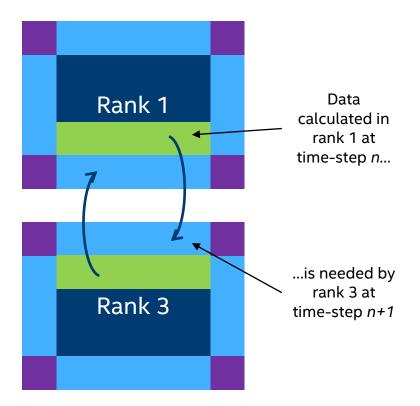
This data changes every time-step and must be exchanged between ranks every time-step to keep the data consistent.

# Schematic of 2D, 2×2 rank y-edge halo exchanges

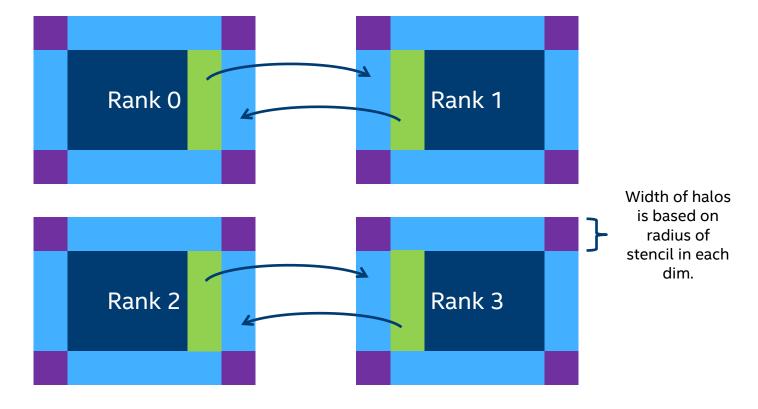
Halos are exchanged at beginning of run and after each time-step.

Can get more complex when there are multiple dependent stencils.



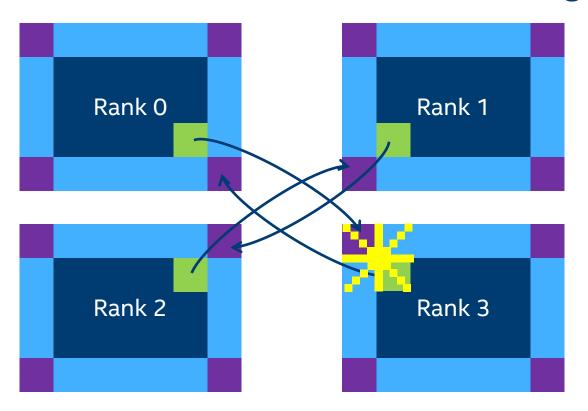


# Schematic of 2D, 2×2 rank x-edge halo exchanges



# Schematic of 2D, 2×2 rank corner exchanges

Corner
exchanges are
only needed
for halos that
have diagonal
points.
This is
determined
automatically
by the stencil
compiler



For 3D stencils, exchanges may be needed along faces, edges, and corners of the rectangular solids.

This concept can be extended to any number of dimensions.

# Running on multiple nodes in a cluster

#### Running the test utility with an explicit MPI command

- bin/yask.sh -mpi cmd 'mpirun <...> -ppn 2' -stencil my stencil -q 1024
  - Replace <...> with the specification of host-names and other needed parameters
    - Use the technique for your scheduler, e.g., -f or -hosts option, or whatever is used on your cluster
    - Important: specify how many processes to run on each node, e.g., with -ppn, usually corresponding to the number of NUMA nodes in each cluster node
  - YASK will automatically partition the global domain and distribute it among the MPI ranks
    - Use the -nr\* YASK options to control the topology of the MPI ranks
    - Example: -nr 2 specifies 2 ranks in each spatial dimension
    - Example: -nrx 2 -nry 4 -nrz 8 specifies 2 ranks in the 'x' dimension, 4 in the 'y', and 8 in the 'z', for 64 total ranks
  - Shared-memory (shm) communication will automatically be applied between ranks that share a virtual address space, but MPI can be forced between all ranks with -no-use shm as explained earlier

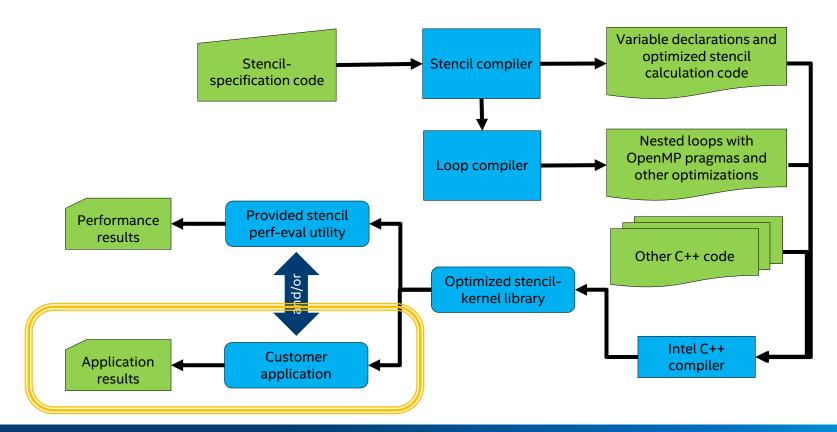
#### Implementation optimization

- For typical configurations, YASK overlaps communication with some computation using this sequence [simplified]:
  - Only the data needed by another rank is calculated on each rank for a given time-step
  - Asynchronous (non-blocking) MPI receive and send requests are initiated
  - The remaining data on each rank is calculated while the MPI data is in-flight
  - The MPI requests are completed before the next time-step is begun
- You can disable this feature via -no-overlap comms to measure the impact



# USING THE APIS

# Making a stencil application



### **API** overview

#### Purpose

- Define the DSL (Domain-Specific Language) used by the YASK compiler
- Facilitate inclusion of stencil code generated by YASK into real applications

#### Target languages

- C++ natively
- Python\* interface generated via SWIG\*

#### Design principles

- Consistent interface
- Stable interface with backward-compatibility maintained except in rare cases
- Documented
- Hidden implementation via C++ pure-virtual classes in most cases
- "Factory" pattern: use small number of factory objects to create more objects



### **API** documentation

From <a href="https://github.com/intel/yask">https://github.com/intel/yask</a>, follow the "API documentation" link in the Overview section of README.md

#### There are four main tabs

- Main Page contains an overview of the YASK workflow and API sets
  - Read this first
- Modules lists the API sets
  - This is the best place to start looking for the API you want
  - YASK Compiler APIs are used to define stencil solutions, i.e., the "DSL"
    - Objects in the compiler API start with "yc"
  - YASK Kernel APIs are used to create your own application from a YASK library
    - Objects in the kernel API start with "yk"
  - YASK Common Utilities contains some common utility classes like output streams
    - Objects here start with "yask\_"
- Classes lists all the C++ classes in alphabetical order
- Files lists the C++ header files



### **API** combinations

#### There are two main steps in using a YASK stencil:

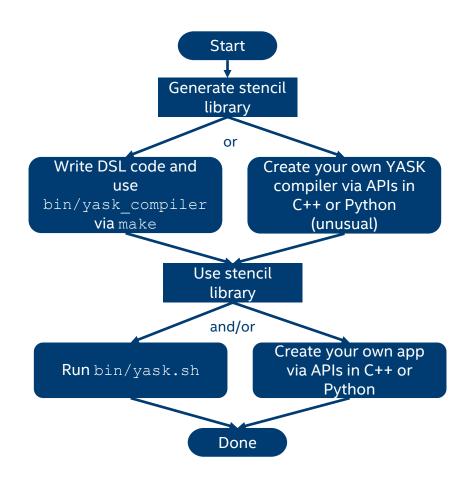
- Create a library from the YASK DSL via the "compiler" APIs
- 2. Use the library in an application via the "kernel" APIs

### For each of these steps, there are three options (making 9 combos):

- Use the binary provided in the YASK package
- Write your own app in C++
- Write your own app in Python

#### The most common combos:

- 1. Compiler: Use the provided YASK compiler binary to create a library for your stencil via the Makefile
- Kernel:
  - Use the provided YASK binary to test and tune your stencil's performance by running yask.sh
  - Create your own binary using the C++ kernel APIs to use the stencil library in production





### Access to the APIs

#### C++

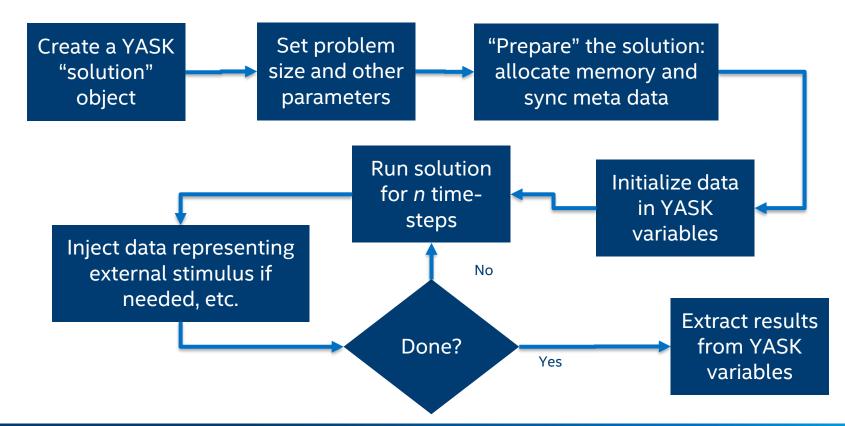
- The standard method of writing YASK stencils uses the "DSL" part of the C++ compiler API library
  - Many examples in the src/stencils directory
  - The provided YASK compiler generates C++ stencil code from the DSL code; this is compiled into a kernel library
  - You only need to make your own YASK compiler binary in special situations, such as creating a custom stencil generator; otherwise, just build your stencil from the Makefile and use the generated library
    - The rest of this section in this tutorial covers using the kernel APIs, not the compiler APIs
- When you run make as in the previous examples, you have implicitly made a kernel library containing the kernel APIs
  - A shared object is created in the lib directory, one for each stencil and target architecture
  - The performance-test binaries we've been running via yask, sh are just wrappers around the library
- The header files are in the include directory
  - #include "yask kernel api.hpp" in your application
  - See src/kernel/tests/yask kernel api test.cpp for an example

#### **Python**

- Build the APIs for Python\* only if you want to make a Python app instead of a C++ app
- Download src code as before: git clone https://github.com/intel/yask
  - See README.md for SWIG\* version requirement
- From yask directory, run make -j api
- Python APIs are not documented separately—use the C++ docs with appropriate type mapping
- Performance-wise, the stencil calculations will be the same speed as in C++, but data access is slower



# Main YASK kernel process steps



# Creating a YASK kernel via the APIs

#### **Bootstrap**

- Declare an object of type yk factory
- Call yk\_factory::new\_env() to create an "environment" object
  - This can be used to provide an existing MPI communicator if desired
- Call yk\_factory::new\_solution() to create a YASK "solution" object
  - The solution object provides methods to configure the problem size, access YASK variables, and run the stencil calculations

#### FAQ: What's with all the output?

- Calling many APIs results in debug output written to a stream
  - By default, this stream is standard output
  - Handy for development, but you probably don't want this in your final application
- Call yk\_solution::set\_debug\_output() to change it to a file or yk\_solution:: disable\_debug\_output() to simply discard it
  - Use a yask\_output\_factory to create new output streams
  - NB: this mechanism was used instead of raw C++ streams to enable Python usage, but you can get access to the C++ stream via yk env::get debug output()->get ostream()



### Key terms related to domain sizes

#### Overall (global) and rank (local) domain sizes

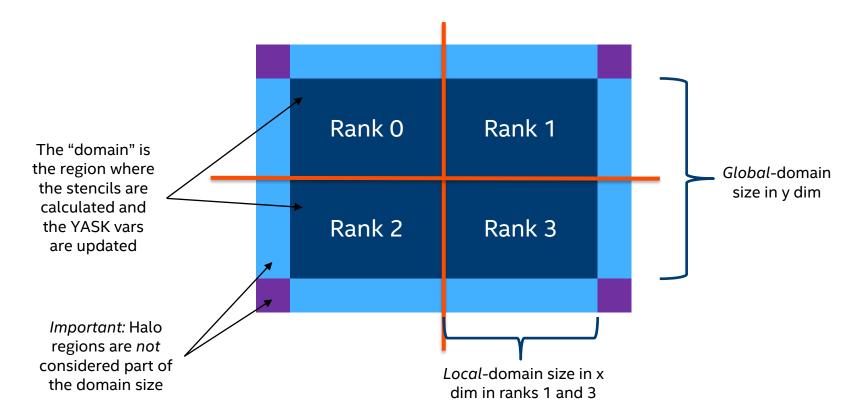
- As with the test utility, each MPI rank can define the local-domain size or the global-domain size in each dimension
  - The size that is not specified will be calculated automatically
  - All ranks in a given row must have the same rank-domain height, etc.
- Also, each rank can specify the number of ranks in each dimension or let the run-time assign them automatically

#### YASK-variable dimensions and sizes

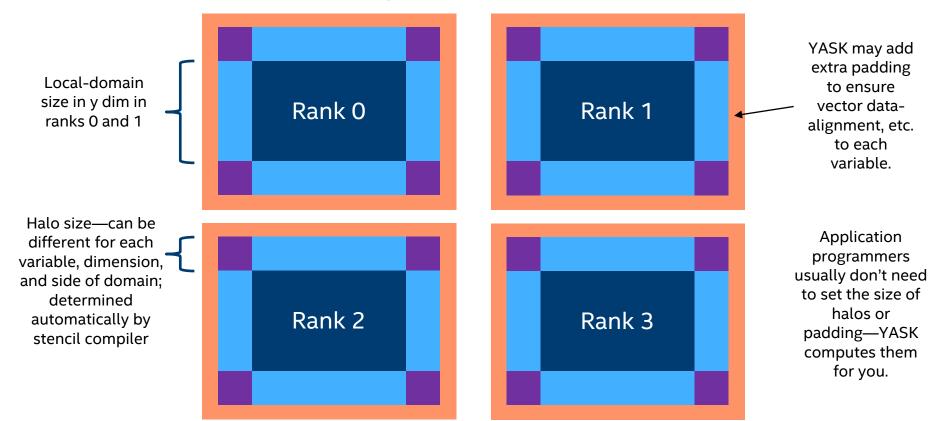
- When YASK variables are defined, they can have three types of dims
  - Step dim: usually "t" for "time"
  - Domain dim: usually used for spatial dims like "x" and "y"
    - Generally, can be any domain-size parameter that may be decomposed over ranks
  - Misc dim: an index that is known at YASK-compile time like a coefficient index
- The size of the domain will be used when allocating each variable
  - Read the "Detailed Description" section of the yk\_var API page for an explanation of each size



# Global view of 2D problem size



# Per-rank view of 2D problem sizes



### Prepare the kernel solution

### Set up the problem sizes and ranks

- Call yk\_solution::set\_overall\_domain\_size() to set the global-domain sizes
  or yk\_solution::set\_rank\_domain\_size() to set the local-domain sizes
- Call yk\_solution::set\_num\_ranks() to explicitly set the number of ranks if desired
- Call yk\_solution::set\_rank\_index() to explicitly set the position of each rank if desired

### Optional: Set other solution parameters

- Call yk\_solution::set\_block\_size() to use the best block size you found during the tuning process
- Alternatively, or in addition, you can use the block-size auto-tuner
  - The auto-tuner can be controlled with APIs
  - There are several ways to use the auto-tuner, discussed in the advanced section
- Other solution parameters, including other tile sizes, can be set via yk solution::apply command line options()



### Allocate data and synchronize info across ranks

```
Call yk solution::prepare solution()
```

- Shares MPI positions across ranks or calculates a default position for each rank
  - Since prepare solution() uses MPI calls, it is critical to call it from each rank
- Ensures domain-size consistency across ranks
- Allocates data for each YASK variable based on halo sizes determined by the YASK compiler and the local domain sizes
- Determines MPI buffer requirements and sizes and allocates data for them

### Initialize data

#### Access YASK variables

- Get a list via yk\_solution::get\_vars() or find a specific var by name via yk\_solution::get\_var()
- Either one returns pointers to variables of type yk\_var
  - Used to find all meta data about each var
  - Use get \* index() APIs to determine valid indices
  - Use idx t as a type for indices: size t and int are not good alternatives.

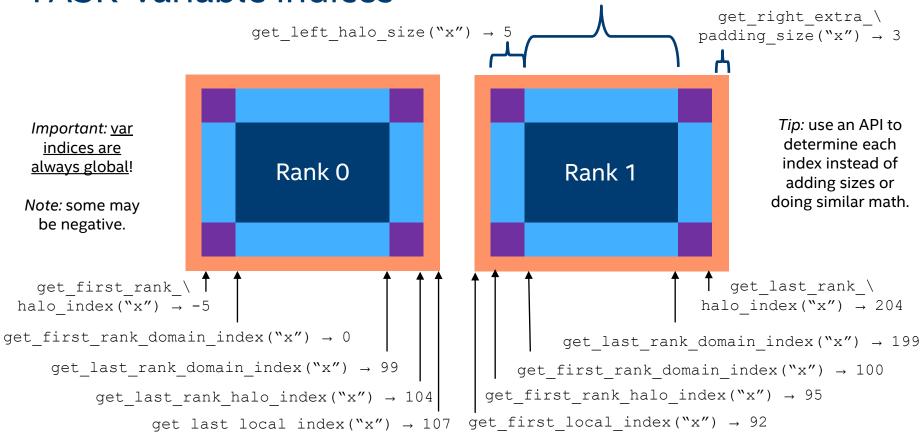
#### Data access

- YASK uses a non-standard tiled data layout ("vector folding" discussed later)
  - Thus, it does not support simple overlay with native row-major or column-major arrays of floating-point numbers
  - NB, github issue #147 requests to support this under certain restrictions for easier integration with legacy applications
- There are several APIs for writing one or more FP elements from and to the variables
  - Start with set\_element(), set\_elements\_in\_slice(), and set\_all\_elements\_same() in yk\_var
  - The "slice" versions are threaded for performance and convert to and from row-major layout
    - But you'll probably want to avoid copying whole vars this way to avoid wasting memory

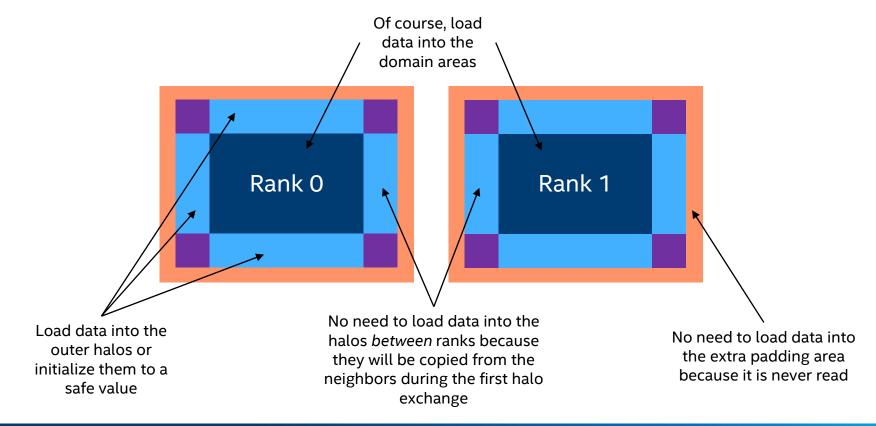


### YASK-variable indices

get rank domain size("x")  $\rightarrow$  100



# Notes on initializing data



### Make the stencil calculations and get results

### Advance the simulation through a series of time-steps

- Call yk solution::run solution()
  - General form takes first and last time indices, e.g.,
    - run solution(0, 9) runs the first 10 time-steps
    - run solution(10, 19) runs the next 10
- If needed, between time-steps, you can access var data, e.g., for source injection

#### Access results

- Use appropriate var methods analogous to the data-writing ones listed earlier
  - Start with get\_element() and get\_elements\_in\_slice()
- Call yk\_solution::get\_stats() to collect some performance stats if desired
- Call yk solution::end solution() when all done to release memory, etc.



### **Advanced APIs**

### Misc interesting APIs (not exhaustive)

- Call yk\_var::set\_numa\_preferred() to set the NUMA node for a given variable—useful for explicit placement in MCDRAM or DDR on Xeon Phi CPUs
- Call yk\_solution::new\_var() to make variables beyond those used in the solution
- Call yk\_solution::apply\_command\_line\_options() to parse a command-line string—useful for quickly applying options from the test utility in another application or for setting options that do not have corresponding APIs
- Call yk\_var::add\_to\_element() to atomically update a var element—useful for threaded source injection

### Exceptions

### YASK APIs that can trigger errors throw exceptions

- Allows applications to catch and process errors
- Find exception documentation in the YASK Common Utilities module tab

#### C++

- Throws object of type yask\_exception
- Call yask exception::get message() for a human-readable explanation
- See src/kernel/tests/yask kernel api exception test.cpp

### Python

- Throws RuntimeError as defined by SWIG\*
- Call format() for a human-readable explanation
- See src/kernel/tests/yask kernel api exception test.py



# ADVANCED STENCILS AND TUNING

### Solutions with multiple stencils

#### Purpose

Many modern seismic simulations require multiple variables to be updated, each with a different stencil

#### Independent stencils

- Definition: for any two stencils, the input of one stencil does not depend on the output of the other within the same time-step, and vice-versa
- Example: x-stress(t+1) depends on x-stress(t), and y-stress(t+1) depends on y-stress(t), but x-stress(t+1) does not depend on y-stress(t+1) or vice-versa
  - See define str TL() in src/stencils/SSGElasticStencil.cpp
  - Thanks to contributor Albert Farres from the Barcelona Supercomputing Center!
- For independent stencils, YASK applies the stencils simultaneously in one pass over the domain

#### Dependent stencils

- Definition: for any two stencils, the output of one stencil is required for the input of the other stencil in the same time-step (the opposite cannot also be true)
- Example: x-velocity(t+1) depends on x-stress(t), and x-stress(t+1) depends on x-velocity(t+1) —see same DSL file
- For dependent stencils, YASK cannot apply the stencils simultaneously and makes subsequent passes over the domain
- YASK automatically determines the dependencies between equations in the DSL and schedules their stencils



# Absorbing boundaries

#### Used in wave-propagation simulation

- Simulation domain covers only a finite block of earth, water, etc.
- Simulated waves will erroneously reflect from the arbitrary boundaries
- Artificial absorbing boundary conditions (ABC) or boundary layers are used to reduce these reflections
- There are many techniques and papers covering this field of study

#### Two high-level approaches to use in YASK

- Modify the stencil(s) to include simple attenuation factors
  - Example: Cerjan sponge layers
  - See Iso3dfdSpongeStencil class in src/stencils/Iso3dfdStencil.cpp
    - Try modifying your test stencil to use sponge layers and check the impact on performance
  - This stencil illustrates the use of combining 3D and 2D variables to save memory
  - Tip: this stencil and its parent class also show the recommended practice of creating functions to return expressions, allowing more flexible composition
- Use different stencils in the boundary layers by creating sub-domains
  - Examples: Higdon BCs, perfectly-matched-layers (PML), and the convolutional form (CPML)





### Sub-domain (spatial) conditions

A sub-domain is a subset of the domain in which a stencil is applied

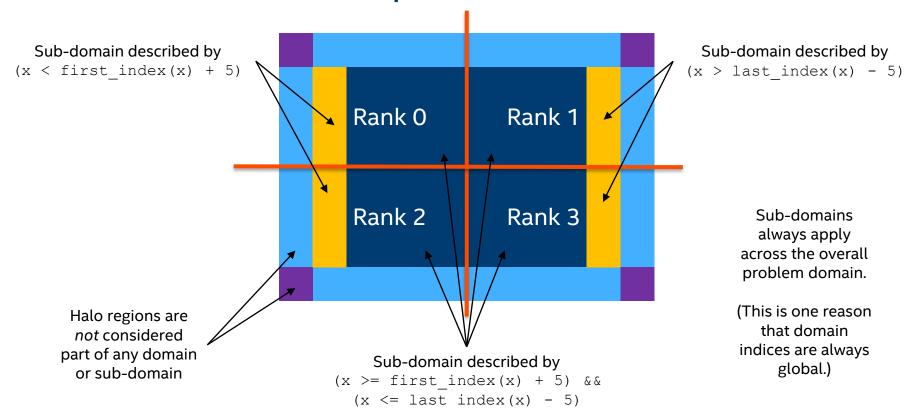
Defining stencils in boundaries is a common use-case for sub-domains

A sub-domain is defined in the DSL by a Boolean expression on the domain indices

- The DSL includes terms for the left-most and right-most indices in the domain
- Example:  $(x < first_index(x) + 5)$  defines a sub-domain on the left side of the domain, 5 elements wide
  - (x > last index(x) 5) would be the same on the right side
  - (x >= first\_index(x) + 5) && (x <= last\_index(x) 5) would be used for the interior sub-domain between the left and right boundary sub-domains
- These conditions are written after the equation using the IF\_DOMAIN operator in the DSL
  - Search for TestBoundaryStencil in src/stencils/TestStencils.cpp for simple synthetic examples
  - See src/stencils/FSGElasticStencil.cpp for a more complex real-world example



### Sub-domains on multiple ranks



# Step (temporal) conditions

### A step-condition expresses which time-steps are valid for a stencil

- Some stencil applications require special processing on regular intervals
- Wave-field sub-sampling can be implemented with step conditions

### Like sub-domain conditions, step conditions restrict stencil application

- Sub-domain expressions can use only the domain indices
- Step-condition expressions can use the step index (usually time) or values in other YASK variables (but not via domain indices)
- Example: (t % 8 == 0) is true every 8<sup>th</sup> time-step
- These conditions are written after the equation using the IF\_STEP operator in the DSL
  - Look for TestStepCondStencil in src/stencils/SimpleTestStencils.cpp for a simple example

# Vector folding (multi-dimensional vectorization)

#### Concept

- Store small 2D or 3D block of data into each SIMD vector
- Pros: reduces memory loads and memory streams compared to traditional 1D in-line vectorization
- Cons: requires non-traditional tiled data layout and additional shift and/or permute operations preceding SIMD arithmetic operations

#### Results

- Significant speedup shown on many systems, esp. those containing high-bandwidth memory (HBM)
- Works well paired with tiling and other performance techniques

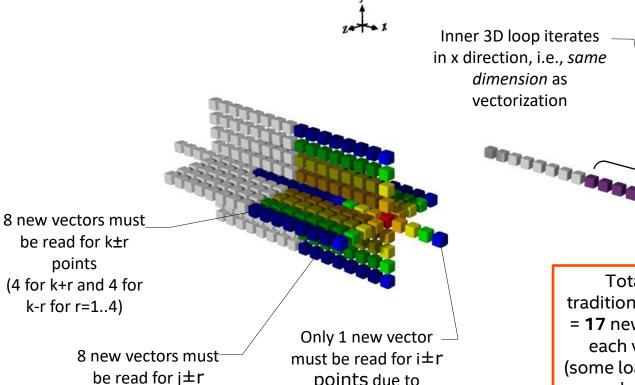
#### **Implementation**

- The YASK compiler automatically generates the proper shift and permute instructions
- The YASK kernel code automatically generates code to store the tiled data and look up elements by index when needed
- Intel® AVX-512 instruction set is needed for efficient permutes, so only enabled when available
- See the paper in the upcoming reading list showing up to  $1.5 \times$  speedup from vector-folding



### **Traditional 1D Vectorization**

points

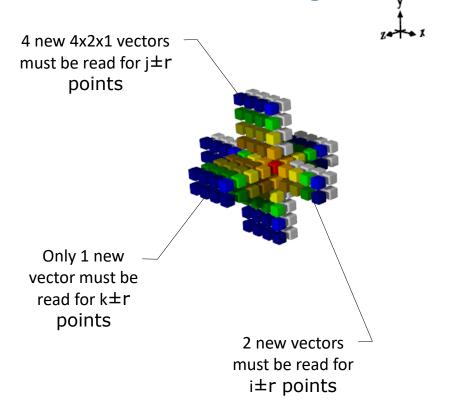


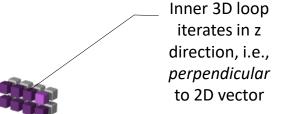
Total BW cost for traditional "in-line" vectors = **17** new vector inputs for each vector of output (some loads will come from cache with blocking)

points due to

overlap along x axis

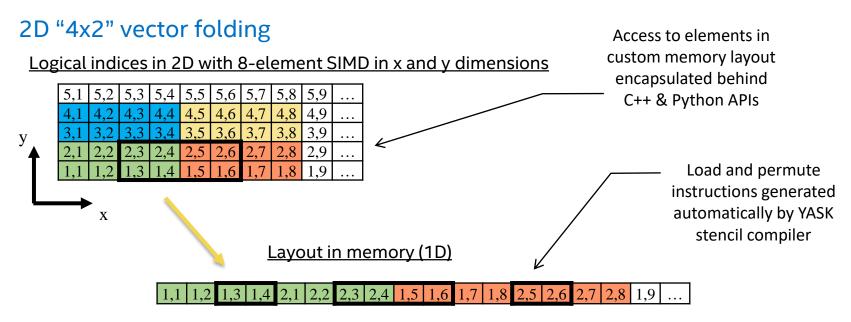
# 2D Vector-Folding





Total BW cost for 4x2x1 vector with z-axis loop= 7 new vector inputs for each vector of output (2.4x lower than in-line)

# Vector-Folding Memory Layout and Code Gen



- 2D vector folding layout (8×1)
- Two aligned vectors are colored
- <u>Unaligned</u> read shown with bold borders done by loading aligned vectors and then shuffling the requisite elements via an AVX-512 permute instruction

### Vector-folding customization

### Specify the vectorization length in each dimension

- Use the fold='x=n, y=n, z=n' argument to the make command-line
  - The values are passed to the YASK stencil compiler and used to generate code
  - The fold settings are also included in non-generated code during compilation
  - Example: make fold='x=1, y=2, z=8' generates code using a 1×2×8 fold
  - Try different fold settings on the test stencil and check the impact on performance
    - Important: Be sure to run make clean before re-compiling when changing compile-time options like vectorization
- The product of the fold lengths should equal the number of SIMD vector elements in the target architecture and FP precision (single or double)
  - Example: single-precision FP using 512-bit SIMD contains 16 elements per vector
  - If the fold is not the right size, the compiler will adjust the requested fold using a heuristic algorithm (i.e., the fold you specify is a hint)
  - The vector length in any dimension not specified defaults to one (1)
  - The default fold varies depending on dimensionality, architecture, and FP precision
    - Is 4×4×1 for a 3D problem using SP FP on Xeon Scalable CPUs, for example





# More compile-time settings

#### Floating-point precision

- The default FP size is 4 bytes, which is the norm in seismic modeling
- Build with double-precision by adding real bytes=8 to the make command

#### Prefetching

- Software prefetch instructions can be added by the YASK compiler
  - Use make ... pfd l1=n to set the L1 prefetch distance to n iterations ahead
  - Use make ... pfd\_12=n to set the L2 prefetch distance to n iterations ahead
  - Set n to zero (0) to disable generation of any prefetch instructions for a level
- The default prefetch depends on the architecture

#### Time allocation

- By default, the YASK compiler tries to determine the minimum number of time-steps that need to be saved in each variable, but you may want to override it
- Override this with make ... time\_alloc=n (to affect all vars at compile-time)
- Override at run-time with the yk\_var::set\_alloc\_size() API



# More compile-time settings

#### Order of domain dimensions

- The order of domain dimensions determines
  - The inner-most or "unit-stride" dimension in the memory layout
  - The nesting order of loops for the stencil-calculation code
  - Default vector-folding and default rank layout
- By default, the order of domain dimensions is determined from the variable declarations
  - For example, new var ("A", {t, x, y, z}) implies domain dimensions are in 'x, y, z' order
  - If vars have different domain-dimension index orders, dimensions are sorted in the order they are seen by the compiler
- You can explicitly set the order of domain dimensions
  - For example, make ... domain dims='z,y,x' creates domain dimensions in 'z,y,x' order
  - This setting overrides the order of the dimensions specified in the variable declarations and affects memory layout, looping order, vector-folding, and rank layout
  - Try different orders on the test stencil and check the performance impact (remember to make clean)

#### Misc settings

- Any preprocessor macro can be set by make ... EXTRA MACROS='name=value ...'
- Type make help for examples of settings the C++ optimization level and more



# More grouping terms: clusters, bundles, and stages

#### **Clusters**

- The YASK compiler can also unroll the innermost loops a specified number of times, effectively calculating a cluster
  of SIMD vectors in each iteration
- Control via the cluster='x=n, y=n, z=n' argument to the make command-line similar to the fold argument for vector-folding
  - Try different cluster settings on the test stencil and check the performance impact (remember to make clean)
- The default clustering is to not unroll (cluster size of one in each dimension)
- Clustering increases register pressure, making it useful only for small stencils

#### **Bundles**

- A bundle is the term used for independent equations that are grouped into one function by the YASK compiler
- A bundle cannot contain equations from different sub-domains
- The compiler eliminates common sub-expressions between stencils in a bundle

#### Stages

- A stage is the term used for a group of equation bundles that are not inter-dependent
- This can occur when there are equations in different sub-domains that are not dependent on one another
- Two or more stages may be evaluated within one cache block that crosses sub-domain boundaries
- Multiple stages are scheduled in the correct order based on dependencies (e.g., in a staggered-grid problem)



# Temporal wave-front rank tiling

#### Goal

- Increase temporal locality in caches at a package level
- Especially useful for large unified cache (such as HBM used for caching DDR memory)

### Technique

- Evaluate n time-steps within a subset of the local rank domain
  - We refer to this subset as a Mega-block in YASK
- Evaluate each Mega-block sequentially until entire rank domain is evaluated for n timesteps

### Increased software complexity

- Must respect mathematical dependencies between time-steps
- Halo exchanges between ranks become more complex
- Any inter-time-step data processing (like source injection) becomes more complex

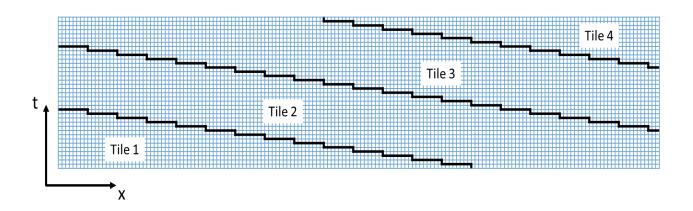


#### Temporal wave-front dependencies Key o value already known value not yet known Highest x-value that may be known value used as input calculated for t=4 due to the value being calculated dependencies on t=3 values 0000000 t=4 00000000000000000 t=3 t=2 Highest x-value that may be calculated for t=2 due to the dependencies on t=1 values Tile ends here for t=1

- In the first temporal wave-front tile shown here, the number of values that can be calculated is reduced for each time-step (and/or each stage within a time-step)
- The amount of shift is called the *wave-front or temporal angle* and is based on the radius of the largest stencil (shift of 8 shown)



# Covering a temporal range via multiple tiles



- Wave-front tiles (Mega-blocks) are computed *sequentially*, but multiple values within one time-step of a given tile may still be evaluated *concurrently*
- After last tile is complete, YASK variables contain the same data as they would have without temporal tiling, so no change is needed when inputting data or reading results
- This concept is directly extended to 2D or 3D stencils by shifting in each spatial dimension
- To handle MPI halo exchanges, some redundant calculations are made between ranks; this is handled by YASK automatically

# Using temporal wave-front rank tiling in YASK

### Another level of tiling hierarchy

- Earlier, we explained how each rank domain is divided into *blocks*, set via -b
- Ranks may also be divided into Mega-blocks, where each Mega-block is a wave-front tile
- Control the spatial size of a Mega-block with -Mb, -Mbx, -Mby, etc.
  - Be sure to use a capital "M"; a lower-case "m" is for micro-blocks, explained later
- Control the temporal size of a Mega-block with -Mbt (assuming your time dim is "t")
- By default, the spatial size of a Mega-block is the size of the rank

### Selecting values

- When using a very large unified cache (e.g., HBM used to cache DDR memory), set the spatial size to fit within this cache
  - See the paper in the upcoming reading list showing ~2× speedup on two stencils
- For processors that have much smaller third-level caches (e.g., a Xeon Scalable processor), values are much more critical, and additional performance is more difficult to obtain

# Temporal block-level tiling

#### Goal

• Increase temporal locality in caches at a core level, e.g., level-2 caches

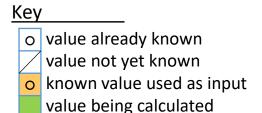
### Technique

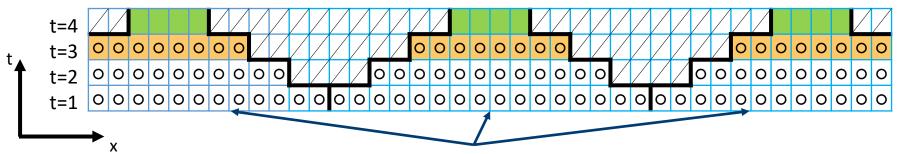
- Evaluate *n* time-steps in each block
  - Recall that a rank is composed of Mega-blocks, and a Mega-block is composed of blocks
- Evaluate blocks concurrently until entire Mega-block is evaluated for n time-steps

### Increased software complexity (similar to wave-front tiling)

- Must respect mathematical dependencies between time-steps
  - Due to concurrency, shapes of blocks are more complex than with sequential wavefront tiling
- Any inter-time-step data processing (like source injection) becomes more complex

# Temporal block dependencies





These three "upward" triangle blocks may be evaluated concurrently. When they are complete, the "downward" triangle blocks are evaluated.

- In one spatial dimension (as shown), this is called triangle or trapezoid tiling (half of diamond tiling)
- Extension into multiple spatial dimensions uses a more complex series of multi-dimensional shapes (polytopes) to tessellate the space (not shown)
- The amount of shift or temporal angle is based on the radius of the largest stencil (shift of 2 shown)

# Using temporal block tiling in YASK

### Not another level of tiling hierarchy

- As before, control the spatial size of a block with -b, -bx, -by, etc.
- Control the temporal size of a block with -bt (assuming your time dim is "t")
- By default, there is no temporal block tiling
- Temporal block tiling occurs within wave-front rank tiling, so the temporal size of a Mega-block defaults to the temporal size of a block
  - But the default spatial size of a Mega-block is still the spatial size of a rank domain
  - You can use both wave-front rank tiling and temporal block tiling, and the temporal size of the Mega-blocks can be larger than the temporal size of the blocks

### Selecting values

 When selecting spatial sizes, consider level-2 cache size, number of YASK variables accessed, the FP-element size, and the number of cores

# Temporal wave-front micro-block tiling

#### Goal

- Observation: block sizes are used as a unit-of-work for OpenMP\* threads as well as level-2 cache targeting
  - These objectives can conflict when there are many variables in a stencil, which requires small blocks to fit in cache
  - This can lead to many short OpenMP tasks, which may not be optimal
- Also, it may be beneficial in some cases to combine the concepts of trapezoid tiling and wave-front tiling across different dimensions
- Goal is to separate the thread and cache block concepts and provide more tiling flexibility

#### Technique

- Evaluate n time-steps in a subset of each block
  - We refer to this subset as a micro-block in YASK
- Evaluate each micro-block sequentially until entire block is evaluated

#### Increased software complexity (similar to previous tiling)

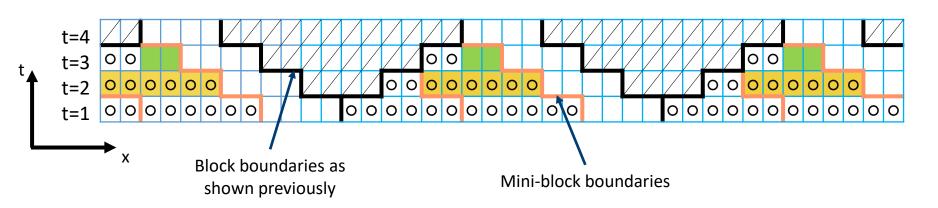
- Must respect mathematical dependencies between time-steps
  - Similar to sequential wave-front tiling, but within [hyper] trapezoids of temporal blocks



# Temporal micro-block dependencies



- o value already known value not yet known
- o known value used as input value being calculated



- Blocks are evaluated *concurrently* as before using OpenMP threads
- Micro-blocks are evaluated sequentially within each block
- Micro-blocks use wave-front tiling, similar to Mega-blocks, but inside blocks instead of ranks



# Using micro-block tiling in YASK

## Another level of tiling hierarchy

- Blocks may be divided into micro-blocks, where each micro-block is a wave-front tile
- Control the spatial size of a micro-block with -mb, -mbx, -mby, etc.
- The temporal size of a micro-block is always the same as a block, so it is set implicitly via -bt
- By default, the spatial size of a micro-block is the size of a block

## Selecting values

- The size of a micro-block should normally correspond to the size of a level-2 cache, considering the number of variables accessed and the FP-element size
  - Any scratch-var values are evaluated at the micro-block level, so consider their sizes
- The block sizes can now be larger than what would fit in a level-2 cache, considering the number of cores with thread balancing
- By sizing micro-blocks as wide as blocks in one or two dimensions and/or sizing blocks as wide as Mega-blocks in some dimensions, interesting special-case scenarios may be created



# Nested OpenMP threads

#### Goal

- Observation: using hyper-threads (SMT) across blocks effectively reduces the usable size of the level-2 cache available to each thread
- This is particularly impactful on Xeon Phi processors with not only 4 hyper-threads per core but also 2 cores sharing a level-2 cache
- Goal: allow multiple threads to use shared caches constructively rather than destructively

## **Technique**

- Evaluate subsets of each micro-block by threads that share caches
  - We refer to this subset as a nano-block in YASK
- Evaluate nano-blocks concurrently until a <u>single time-step</u> in a mini-block is evaluated (thus, there is no temporal nano-block tiling)
- Since blocking already uses OpenMP threads, nano-blocking is implemented with a nested level of OpenMP threading



# Using nano-block tiling in YASK

#### Another level of tiling hierarchy

- Micro-blocks may be divided into nano-blocks
- Control the spatial size of a nano-block with -nb, -nbx, -nby, etc.
- There is no temporal nano-blocking, so there is not an -nt option

#### Selecting values

- By default, the spatial size of a nano-block depends on the number of threads are used per block
  - If there is one thread per block (no nested OpenMP), the default size of a nano-block is the size of a micro-block
    - In this case, it is recommended to use only one thread per level-2 cache
  - If there are >1 threads per block (nested OpenMP active), the default size of a nano-block is a narrow slab the width of
    one vector (usually in the first dim) and the size of a micro-block in the other dims
    - This setting is intended to increase reuse between hyper-threads while keeping threading overhead as low as possible
- Control the number of threads
  - By default, the yask.sh script runs one thread per core by passing -outer\_threads to the binary, and the default number of inner threads is one (1).
  - If you want to use hyper-threads, use the -inner\_threads option, e.g., -inner\_threads 2 will use both hyper-threads on most Intel® Xeon™ processors, assuming hyper-threading is enabled in the BIOS settings
  - You can also override the -outer\_threads value to, for example, experiment with multiple cores sharing data across their L2 caches
- Try adjusting threads-per-block and/or nano-block sizes on the test stencil and check performance





# Using pico-block tiling in YASK

## Lowest level of tiling hierarchy

- Nano-blocks may be divided into pico-blocks
- Control the spatial size of a pico-block with -pb, -pbx, -pby, etc.
- There is no temporal pico-blocking, so there is not an -pt option
- Pico-blocking can be used for level-one (L1) cache-blocking on CPUs, but it's use is more critical on GPUs, explained later in this tutorial

## Selecting values

- By default, there is only one pico-block per nano-block
- Pico-blocking allows each nano-block, which is evaluated by an inner OpenMP thread, to have even more locality
- In some stencils, pico-blocking may be able to take advantage of L1 cache locality
- Try adjusting pico-block sizes on the test stencil and check performance





# Review of the CPU tiling hierarchy

Hierarchy level	Spatial size options	Concurrent evaluation?	Temporal tiling?	Temporal size options (assuming time dim is "t")
Global domain	-g*	N/A (only one)	N/A	-trial_steps
Local domains	-[*	Yes, via MPI	No	
Mega-blocks	-Mb*	No	Yes, via sequential wave- front tiling	-Mbt
Blocks	-b*	Yes, via outer OpenMP threads	Yes, via concurrent hyper- trapezoid tiling	-bt
Micro-blocks	-mb*	No	Yes, via sequential wave- front tiling	-bt (implicitly)
Nano-blocks	-nb*	Yes, via inner OpenMP threads	No	
Pico-blocks	-pb*	No (unless on GPU)	No	
Clusters	cluster=	No	No	
Vectors	fold=	No	No	
Elements	real_bytes=	Yes, via SIMD	No	

# Example tile settings for temporal tiling

The following options were hand-tuned for an Intel® Xeon® Platinum 8352Y CPU

```
bin/yask.sh -stencil my_stencil -1 1024 \
-mbx 32 -mby 32 -mbz 128 \
-b 256 -bt 12 \
-no-pre_auto_tune
```



- Micro-block settings selected targeting 1MiB
  - The Intel® 8352Y has a 1280KiB L2 cache per core, and this gives some extra space
  - 32 × 32 × 128 × 4B × 2 = 1MiB (128 ⇒ longer in unit stride; "4B" ⇒ 4 bytes per FP element; "× 2" to keep two time-steps in cache)
- Spatial block settings selected for a multiple of number of cores per socket
  - The Intel® 8352Y has 32 cores per socket
  - Trying 2 blocks per core: 64 blocks (modify for CPU with different core count)
    - A small multiple (like 2 or 4) often works better than the exact count because several blocks per core gives more opportunity for dynamic load balancing
  - Trying 4 blocks across each dimension: 4 × 4 × 4 = 64
  - 1024 ÷ 4 = 256, so using -b 256
- Temporal block setting (-bt 12) was chosen experimentally by iterating manually through several values
- Mega-block size not set because L3 blocking is not being targeted in this experiment
- Nano-block size not set because only one thread per block is used in this experiment
- Pico-block size not set because L1 blocking is not being targeted in this experiment
- These settings resulted in about a 40% speedup over default non-temporal tiling; your results may vary

# Controlling when the automatic tuner runs

#### Pre-calculation mode

- Runs before the desired stencil calculations are done
- Does not ensure that calculations are done in the proper order
  - Thus, important to [re]initialize data after running in this mode for use in deployed applications
- Intended for benchmarking to tune block-size before running actual time-steps
- Use [-no]-pre auto tune option to control in provided test utility (default is on)
- Call API run auto tuner now() to activate outside in your application

#### Intra-calculation mode

- Runs during desired stencil calculations
- Maintains proper calculations
- Intended for deployment, esp. on multiple platforms or when all final platforms are not known α
  priori
- Use [-no] -auto tune option to control in provided test utility (default is off)
- Call reset\_auto\_tuner() to turn off or on explicitly
- Call API is \_auto\_tuner\_enabled() to determine whether it is [still] running



# Controlling what the automatic tuner does

## Tiling levels

- By default, the auto-tuner is applied to the "block" tiles only
- More generally, you can supply a list of tiling levels via -auto\_tune\_targets
  - The argument is a comma-separated list of level abbreviations: Mb for mega-blocks, b for blocks, mb for micro-blocks, nb for nano-blocks, and pb for pico-blocks
  - The order specified is the order in which the tile sizes are tuned, and targets can be repeated
  - Example: -auto\_tune\_targets b, Mb, b will tune blocks, then mega-blocks, then blocks again

#### Other controls

- -auto\_tune\_radius controls how far the tuner will start exploring from the initial settings
- -auto\_tune\_trial\_secs controls how long a new trial must be run to consider it better than the current best trial; higher values avoid measurement noise but require more time to converge



# Multi-parameter auto-tuner script

#### Goal

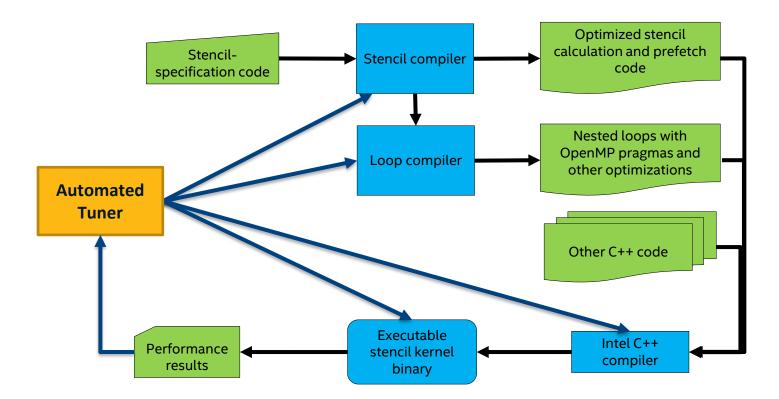
- Use a genetic algorithm to explore the design space of many parameters
- Can target compile-time and/or run-time parameters

## Usage

- Run utils/bin/yask tuner.pl to see options
- By default, explores almost every compile-time and run-time parameter
  - Practically, you probably want to limit some of them
  - To fix the problem size, use -g\* or -1\* options, e.g., -g=1024 -gz=128
  - To search only run-time parameters, use -noBuild option
    - Must build binary before running tuner
- Tuner will stop when there has been no improvement over 5 generations
- Tip: alternate between manual and automated tuning for final optimization
- Currently only works for 3D (spatial) problems with dims named t, x, y, and z



# YASK workflow driven by auto-tuner



# **GPU OFFLOADING**

## **GPU** offload overview

#### Mechanism

- GPU offloading via OpenMP Device Directives (sec. 2.14 in OpenMP 5.1 spec)
  - Build with offload=1 to enable GPU offloading with explicit data transfers
    - Or, build with offload usm=1 to enable offloading with a unified shared memory model
  - Add offload arch=device name to make command to change the OpenMP target
    - Default device name is spir64 to target all Intel GPUs
    - When targeting spir64, specific device code is generated at run-time with JIT offload compiler
- The nano- and pico-block loops are offloaded
  - Nano loops executed via omp target teams distribute
    - Evaluates pico-blocks in parallel by OpenMP target teams
    - Thread limit set via -device\_thread\_limit option
  - Pico loops (nested within nano loops) executed via omp parallel for
    - Pico-block sizes affect GPU-cache locality and data reuse
      - These are the default auto-tuner targets for an offload binary
    - For an N-spatial-dimensional solution, N-1 dim loops are concurrent; one is sequential
  - All higher-level loops are executed on the CPU
    - The default number of CPUs threads for an offload binary is one (1)
    - Using more than one CPU thread allows launching more than one GPU kernel concurrently



# GPU usage examples

#### Build

- make -j stencil=iso3dfd offload=1
  - Note the arch string that is printed at the end of the build

#### Run

- You must specify the arch string for the offload binary
  - The arch string will be the CPU target plus .offload-device\_name; see examples below
  - The default *arch* will select the current CPU target only
- bin/yask.sh -stencil iso3dfd -arch avx512.offload-spir64 ranks 1 -device\_thread\_limit 1024 -g 512 -no-pre\_auto\_tune pbx 256 -pby 32 -pbz 32
- bin/yask.sh -stencil iso3dfd -arch avx512.offload-spir64 ranks 1 -device\_thread\_limit 1024 -g 512 -pre\_auto\_tune

# WRAP-UP

# Read more about YASK features and applications

- "Vector Folding: improving stencil performance via multi-dimensional SIMD-vector representation."
   C Yount. 17th International Conference on High Performance Computing and Communications (HPCC), 2015
- "YASK—Yet Another Stencil Kernel: A Framework for HPC Stencil Code-Generation and Tuning." C Yount, J Tobin, A Breuer, A Duran. Domain-Specific Languages and High-Level Frameworks for High Performance, 2016
- "Effective use of large high-bandwidth memory caches in HPC stencil computation via temporal wave-front tiling." C Yount, A Duran. 7th International Workshop on Performance Modeling, Benchmarking and Simulation of High Performance, 2016
- "Accelerating seismic simulations using the Intel Xeon Phi knights landing processor." J Tobin, A Breuer, A Heinecke, C Yount, Y Cui. International Supercomputing Conference, 2017
- "Performance Optimization of Fully Anisotropic Elastic Wave Propagation on 2nd Generation Intel® Xeon Phi (TM) Processors." A Farres, C Rosas, M Hanzich, A Duran, C Yount. 2018 IEEE International Parallel and Distributed Processing Symposium
- "Multi-level spatial and temporal tiling for efficient HPC stencil computation on many-core processors with large shared caches." C Yount, A Duran, J Tobin. Future Generation Computer Systems, March, 2019



## Call to action

## Work through this tutorial

## Code your own stencil

- Use it in a real application
- Please feel free to contact the developers with questions
- Contribute your stencil for others to use: create a fork on github\* and submit a pull request
- Please tell the developers about your experience, maybe even co-publish results

## Contribute to the project

- See the "issues" database on github\* for the to-do list
- Talk to the developers about something you'd like to work on
  - Probably some good academic projects in there!





https://github.com/intel/yask

