

High Performance Computing: Speed Optimisation Using NumPy, MPI And GPU Acceleration Day 3

Advanced Scientific Programming with Python

Class Inheritance

- Classes can inherit from other classes.
- A class can inherit attributes and behaviour methods from another class, called the superclass or parent class.
- A class which inherits from a superclass is called a subclass, also called heir class or child class.
- We could implement a vehicle class in Python, which might have methods like accelerate and brake. Cars, Buses and Trucks and Bikes can be implemented as subclasses which will inherit these methods from vehicle.

Class Inheritance

```
class Pet(object):
    def __init__(self, name, species):
        self.name = name
        self.species = species
    def getName(self):
        return self.name
    def getSpecies(self):
        return self.species
    def __str__(self):
        return "%s is a %s" % (self.name, self.species)
class Dog(Pet):
    def __init__(self, name, chases_cats):
        Pet.__init__(self, name, "Dog")
        self.chases cats = chases cats
    def chasesCats(self):
        return self.chases cats
class Cat(Pet):
    def __init__(self, name, hates_dogs):
        #Pet.__init__(self, name, "Cat")
        super(Cat, self). init (name, "Cat")
        self.hates_dogs = hates_dogs
    def hatesDogs(self):
        return self.hates dogs
```

Multiple Inheritance

Python also support multiple inheritance

```
class Reader(object):
    def init (self):
        self. size = 0
    @property
    def size(self):
        return self._size
    def read(self):
        data = 'data read'
        return data
class Writer(object):
    def __init__(self):
        self. size = 0
    @property
    def size(self):
        return self. size
    def write(self, data):
        # write data somewhere
        pass
class ReaderWriter(Reader, Writer):
    def __init__(self):
        super(ReaderWriter, self). init ()
```

Why NumPy?

```
import numpy as np

In [4]: %timeit a.sum()
1000 loops, best of 3: 415 μs per loop

a = np.random.random((1000000))
b = list(a)
%timeit a.sum()
%timeit sum(b)

In [5]: %timeit sum(b)
10 loops, best of 3: 41.3 ms per loop
%timeit sum(b)

In [6]: 41.3/0.415
Out[2]: 99.51807228915662
```

How much faster is the NumPy version?



```
A - About 2x
```

B - About 5x

C - About 25x

D - About 100x

E - NumPy is slower

What Is NumPy?

- The numpy package is used in almost all numerical computation using Python.
- It provide high-performance vector, matrix and higher-dimensional data structures for Python.
- It is implemented in C and Fortran so when calculations are vectorized (formulated with vectors and matrices), **performance is very good**.

- To achieve it's high performance NumPy arrays are much more restrictive about data types than Python lists
- All the elements have the same type
- The most common NumPy data types are:
 - numpy.float64, numpy.float32 (also known as double and single precision. The default python float is also 64-bits long)
 - numpy.int64, numpy.uint64, and it's 32 and 16 bit equivalents
 - numpy.complex128 and numpy.complex64 (two float64s and two float32s)
- These are closely related to C types
- The suffix shows the size of the type in bits

- A floating point number x is generally represented as:
 x = sign * mantissa * base**exponent
- They are stored in memory as a sign bit followed by the exponent bits and finally the mantissa or significand. For example for float32:



- The number of bits of the exponent defines the range of the type
- The size of the mantissa defines its precision, or the number of significant digits of the type
- Certain values are values are reserved for a special meaning. The example above typically means NaN (not a number, like 0/0)

- np.uint64 corresponds to an unsigned 64-bit integer
- This is probably the simplest type, using 64bits for the number, it simply encodes integers between 0 and 2**64-1.
- Just be careful with negative numbers or overflow problems, e.g.:

```
In [142]: a = np.array([-1], dtype=np.uint64)
In [143]: a
Out[143]: array([18446744073709551615], dtype=uint64)
In [144]: a += 1
```

- np.uint64 corresponds to an unsigned 64-bit integer
- This is probably the simplest type, using 64bits for the number, it simply encodes integers between 0 and 2**64-1.
- Just be careful with negative numbers or overflow problems, e.g.:

```
In [142]: a = np.array([-1], dtype=np.uint64)
In [143]: a
Out[143]: array([18446744073709551615], dtype=uint64)
In [144]: a += 1
```

```
In [145]: a
Out[145]: array([0], dtype=uint64)
```

- np.int64 corresponds to a signed 64-bit integer
- It encodes integers with 1 bit for the sign and 63 bits for the number, with a range between -2**63 and 2**63-1.
- You can find this with:

```
In [100]: np.iinfo(np.int64)
Out[100]: iinfo(min=-9223372036854775808, max=9223372036854775807, dtype=int64)
```

• The only thing to keep in mind is avoiding overflow problems, e.g.:

```
In [103]: a = np.int64(2**63-1)
In [104]: a
Out[104]: 9223372036854775807
In [105]: a += 1
<ipython-input-105-8076006b0952>:1: RuntimeWarning: overflow encountered in
long_scalars
    a += 1
In [106]: a
Out[106]: -9223372036854775808
```

• This means floating point arithmetic does not always behave like you expect. Adding and subtracting different values can give 0:

```
In [88]: a = np.array([2**24+1],dtype=np.float32)
In [89]: a[0] -= 2**24
In [90]: a[0]
Out[90]: 0.0
```

The result of a sum can depend on the order:

```
In [159]: a = np.array(np.random.rand(100000),dtype=np.float16) - 0.5
In [160]: a.sum()
Out[160]: -86.2
In [161]: np.sort(a).sum()
Out[161]: -93.44
```

Numbers that should be the same are not:

```
In [162]: 1/3 == 1-2/3
Out[162]: False
```

```
In [159]: a = np.array(np.random.rand(100000),dtype=np.float16) - 0.5
In [160]: a.sum()
Out[160]: -86.2
In [161]: np.sort(a).sum()
Out[161]: -93.44
```

Which is the more accurate result sum in the previous example? Why?



A - The unsorted sum

B - The sorted sum

```
In [159]: a = np.array(np.random.rand(100000),dtype=np.float16) - 0.5
In [160]: a.sum()
Out[160]: -86.2
In [161]: np.sort(a).sum()
Out[161]: -93.44
```

Which is the more accurate result sum in the previous example? Why?



A - The unsorted sum

B - The sorted sum

```
In [165]: a.sum(dtype=np.float64)
Out[165]: -86.242919921875

In [166]: np.sort(a).sum(dtype=np.float64)
Out[166]: -86.242919921875
```

- NumPy offers more indexing facilities than regular Python sequences.
- In addition to indexing by integers and slices, arrays can be indexed by arrays of integers and arrays of booleans.

You can also use indexing with arrays as a target to assign to:

```
>>> a = np.arange(5)
>>> a
array([0, 1, 2, 3, 4])
>>> a[[1,3,4]] = 0
>>> a
array([0, 0, 2, 0, 0])
```

- When we index arrays with arrays of (integer) indices we are providing the list of indices to pick.
- With boolean indices the approach is different; we explicitly choose which items in the array we want and which ones we don't.
- The most natural way one can think of for boolean indexing is to use boolean arrays that have the same shape as the original array:

 Fancy indexing is the most general selection method, but it is also the slowest.

```
# Let's create an array with a large number of rows.
# We will select slices of this array along the first dimension.
n, d = 100000, 100
a = np.random.random sample((n, d)); aid = id(a)
# Let's select one every ten rows, using two different methods
# (array view and fancy indexing).
b1 = a[::10]
b2 = a[np.arange(0, n, 10)]
np.array_equal(b1, b2)
True
# Let's compare the performance of both methods.
%timeit a[::10]
1000000 loops, best of 3: 804 ns per loop
%timeit a[np.arange(0, n, 10)]
100 loops, best of 3: 14.1 ms per loop
```

 Fancy indexing is several orders of magnitude slower as it involves copying a large array.

```
>>> a = np.arange(5)
>>> a[[0,0,2]]=[1,2,3]
>>> a
>>> a
>>> a
array([2, 1, 3, 3, 4])
```

```
Socrative UU1

What's the value of a[0]?
```

```
>>> a = np.arange(5)
>>> a[[0,0,2]]+=1
>>> a
array([1, 1, 3, 3, 4])
```



Views

- A view is simply another way of viewing the data of the array.
- The data of both objects are shared.
- You can create views by selecting a slice of the original array or by changing the data type.
- Slice views are the most common.
- The rule of thumb for creating a slice view is that the viewed elements can be addressed with offsets, strides, and counts in the original array.
 For example:

Slice Views

- Are the most common view.
- The viewed elements must be able to be addressed with offsets, strides, and counts in the original array.

```
>>> a = numpy.arange(10)
>>> a
array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
>>> v1 = a[1:2]
>>> v1
array([1])
>>> a[1] = 2
>>> v1
array([2])
>>> v2 = a[1::3]
>>> v2
array([2, 4, 7])
>>> a[7] = 10
>>> v2
array([ 2, 4, 10])
```

Any selection which involves only slices (e.g. 0:10, ::-1, etc...) returns a view.

Data Type Views

- Sometimes you want to inspect memory with a different data type
- This is when data type views come in

You get the raw byte stream casted to whatever data type you selected

Data Ownership

- Now you see that multiple arrays can modify the same data
- Who owns the data?

```
In [20]: b.flags
Out[20]:
  C CONTIGUOUS : True
  F CONTIGUOUS : True
  OWNDATA : True 	
  WRITEABLE : True
  ALIGNED: True
  UPDATEIFCOPY : False
In [21]: v4.flags
Out[21]:
  C CONTIGUOUS : True
  F_CONTIGUOUS : True
  OWNDATA : False ◀
  WRITEABLE: True
  ALIGNED: True
  UPDATEIFCOPY : False
```

View And Data Ownership

```
In [26]: a = numpy.array([1])
In [27]: b = a[:]
In [28]: b[0] = 0
In [29]: a[0] == 0
Out[29]: True
```



What is the result of the code?

```
In [36]: a = numpy.random.random((3))
In [37]: b = a[::-1]
In [38]: b[0] = 1
In [39]: b[0] == a[-1]
Out[39]: True
```

And now?

View And Data Ownership

```
In [40]: a = numpy.random.random((3))
In [41]: b = a
In [42]: b[0] = 1
In [43]: b[0] == a[0]
Out[43]: True
```

```
Socrative UU1

And in this one?
```

```
In [44]: a = numpy.random.random((3))
In [45]: b = a[[0,1,2]]
In [46]: b[0] = 1
In [47]: b[0] == a[0]
Out[47]: False
```

And finally?

Some Tricky Cases

Both snippets on the right seem to do the same thing but they don't.

```
>>> a = numpy.arange(10)
>>> a[[1,2]] = 100
>>> a
array([ 0, 100, 100,  3,  4,  5,  6,  7,  8,  9])

>>> c1 = a[[1,2]]
>>> c1[:] = 100
>>> a
array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
>>> c1
array([100, 100])
```

Some Tricky Cases

Is a[0,0] also 100 in this case?

Broadcasting

- NumPy operations are usually done on pairs of arrays on an elementby-element basis.
- In the simplest case, the two arrays must have exactly the same shape, as in the following example:

```
>>> a = np.array([1.0, 2.0, 3.0])
>>> b = np.array([2.0, 2.0, 2.0])
>>> a * b
array([ 2., 4., 6.])
```

- NumPy's broadcasting rule relaxes this constraint when the arrays' shapes meet certain constraints.
- The simplest broadcasting example occurs when an array and a scalar value are combined in an operation:

```
>>> a = np.array([1.0, 2.0, 3.0])
>>> b = 2.0
>>> a * b
array([ 2., 4., 6.])
```

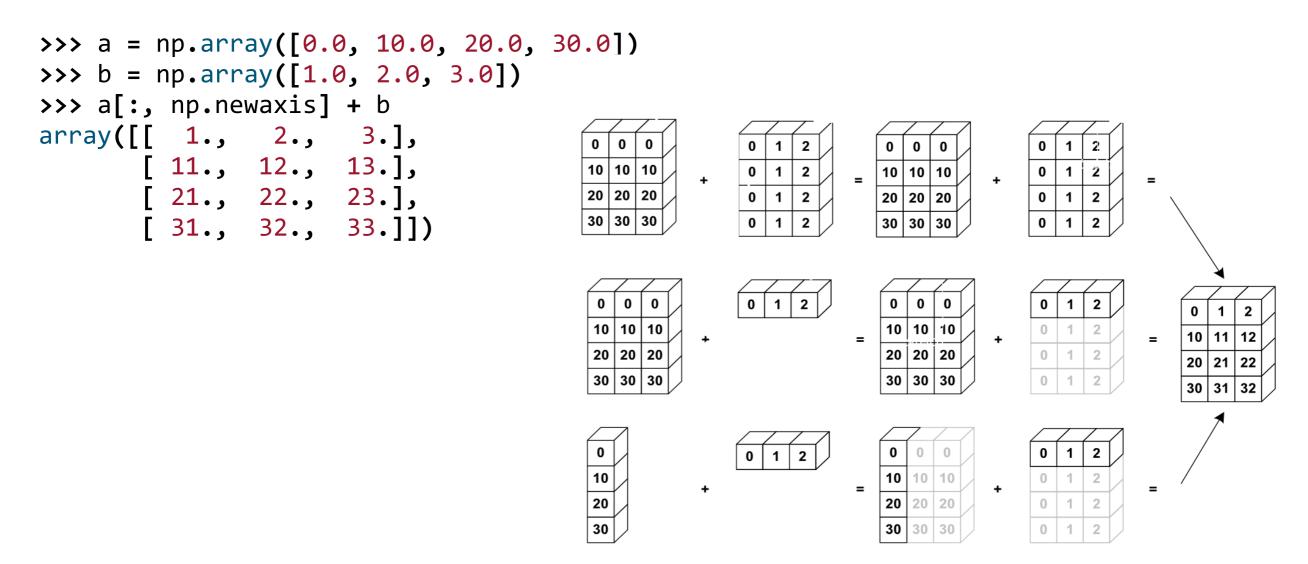
General Broadcasting Rules

- When operating on two arrays, NumPy compares their shapes element-wise.
- It starts with the trailing dimensions, and works its way forward.
- Two dimensions are compatible when:
 - 1. they are equal, or
 - 2. one of them is 1
- The size of the resulting array is the maximum size along each dimension of the input arrays.
- Arrays do not need to have the same *number* of dimensions. For example:

```
Image (3d array): 256 x 256 x 3
Scale (1d array): 3
Result (3d array): 256 x 256 x 3
```

Broadcasting

- Broadcasting provides a convenient way of taking the outer product (or any other outer operation) of two arrays.
- The following example shows an outer addition operation of two 1-d arrays:



Broadcasting

 Here's a comparison between broadcasting and explicit tiling for the outer product:

```
n = 1000
a = np.arange(n)
ac = a[:, np.newaxis]
ar = a[np.newaxis, :]
%timeit np.tile(ac, (1, n)) * np.tile(ar, (n, 1))
100 loops, best of 3: 10 ms per loop
```

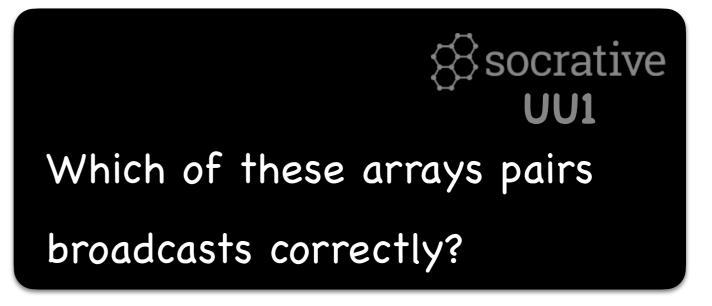
And now using broadcasting

```
%timeit ar * ac
100 loops, best of 3: 2.36 ms per loop
```

It's also much shorter to write!

Broadcasting Exercises

```
(4d array): 8 x 1 x 6 x 1
            (3d array): 7 x 1 x 5
            (2d array): 5 x 4
B
            (1d array):
            (1d array): 3
            (1d array):
            (2d array): 2 x 1
            (3d array): 8 \times 4 \times 3
            (3d array): 15 \times 3 \times 5
E
            (3d array): 15 x 1 x 5
```



Boolean Operations

```
>>> import numpy as np
>>> a = np.array([1, 1, 0, 0], dtype=bool)
>>> b = np.array([1, 0, 1, 0], dtype=bool)
>>> np.logical or(a, b)
array([ True, True, False], dtype=bool)
>>> a + b
array([ True, True, False], dtype=bool)
>>> a | b
array([ True, True, True, False], dtype=bool)
>>> np.logical and(a, b)
array([ True, False, False, False], dtype=bool)
>>> a * b
array([ True, False, False], dtype=bool)
>>> a & b
array([ True, False, False], dtype=bool)
>>> np.logical not(a)
array([False, False, True, True], dtype=bool)
>>> -a
array([False, False, True, True], dtype=bool)
>>> ~a
array([False, False, True, True], dtype=bool)
```

There are many ways to express boolean operation in NumPy.

Avoid Unnecessary Array Copies

The following function can be used to check if two arrays share the same data:

```
def id(x):
    # This function returns the memory
    # block address of an array.
    return x.__array_interface__['data'][0]

a = np.zeros(10);
aid = id(a);

b = a.copy();
id(b) == aid
>>> False
```

Avoid Unnecessary Array Copies

Array computations can involve in-place operations (the array is modified):

```
a *= 2; id(a) == aid
>>> True
```

or implicit-copy operations (a new array is created):

```
c = a * 2; id(c) == aid
>>> False
```

Be sure to choose the type of operation you actually need.

Implicit-copy operations are significantly slower, as shown here:

```
%%timeit a = np.zeros(10000000)
a *= 2
>>> 10 loops, best of 3: 19.2 ms per loop

%%timeit a = np.zeros(10000000)
b = a * 2
>>> 10 loops, best of 3: 42.6 ms per loop
```

How It Works?

- A NumPy array is basically described by metadata (number of dimensions, shape, data type, and so on) and the actual data.
- The data is stored in a homogeneous and contiguous block of memory, at a particular address in system memory (Random Access Memory, or RAM).
- This block of memory is called the data buffer. This is the main difference with a pure Python structure, like a list, where the items are scattered across the system memory.
- This aspect is the critical feature that makes NumPy arrays so efficient.

How It Works?

Why is this so important? Here are the main reasons:

- 1. Array computations can be written very efficiently in a low-level language like C (and a large part of NumPy is actually written in C). Knowing the address of the memory block and the data type, it is just simple arithmetic to loop over all items, for example. There would be a significant overhead to do that in Python with a list.
- 2. Spatial locality in memory access patterns results in significant performance gains, notably thanks to the CPU cache. Indeed, the cache loads bytes in chunks from RAM to the CPU registers. Adjacent items are then loaded very efficiently (sequential locality, or locality of reference).
- 3. Data elements are stored contiguously in memory, so that NumPy can take advantage of vectorized instructions on modern CPUs, like Intel's SSE and AVX, AMD's XOP, and so on. For example, multiple consecutive floating point numbers can be loaded in 128, 256 or 512 bits registers for vectorized arithmetical computations implemented as CPU instructions.

What Is The Difference Between In-Place And Implicit-Copy Operations?

- An expression like a *= 2 corresponds to an in-place operation, where all values of the array are multiplied by two.
- By contrast, a = a * 2 means that a new array containing the values of
 a * 2 is created, and the variable a now points to this new array.
- The old array becomes unreferenced and will be deleted by the garbage collector.
- No memory allocation happens in the first case, contrary to the second case.
- More generally, expressions like a[i:j] are views to parts of an array: they point to the memory buffer containing the data.
- Modifying them with in-place operations changes the original array. Hence, a[:] = a * 2 results in an in-place operation, unlike a = a * 2.

What Is The Difference Between In-Place And Implicit-Copy Operations?

Knowing this subtlety of NumPy can help you fix some bugs (where an array is implicitly and unintentionally modified because of an operation on a view), and optimize the speed and memory consumption of your code by reducing the number of unnecessary copies.

```
import numpy as np
a = np.array([1])
b = np.array([0])
print(a[0]+1.5)
b = a[:]
a += 1
print(a[0])
print(b[0])
```



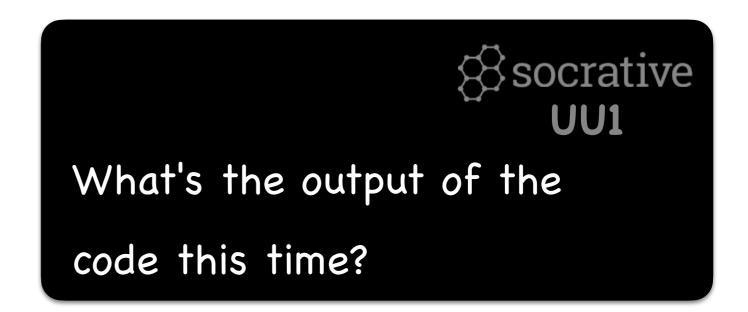
What Is The Difference Between In-Place And Implicit-Copy Operations?

Note also the important difference between a and b in the following lines of code:

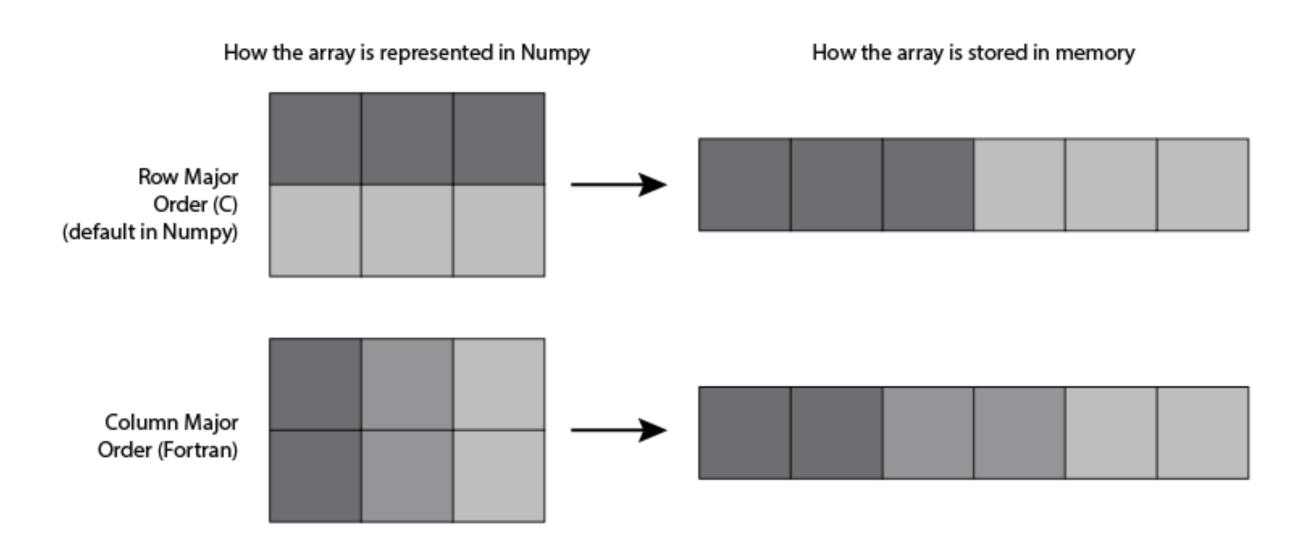
```
import numpy as np
a = np.array([0])
b = np.array([0])

a += 0.5
b = b + 0.5

print(a)
print(b)
```



- A 2D matrix contains items indexed by two numbers (row and column), but it is stored internally as a 1D contiguous block of memory, accessible with a single number.
- There is more than one way of storing matrix items in a 1D block of memory: we can put the elements of the first row first, the second row then, and so on, or the elements of the first column first, the second column then, and so on.
- The first method is called row-major order, whereas the latter is called column-major order.
- Choosing between the two methods is only a matter of internal convention: NumPy uses the row-major order, like C, but unlike FORTRAN.



- More generally, NumPy uses the notion of strides to convert between a multidimensional index and the memory location of the underlying (1D) sequence of elements.
- The specific mapping between array[i1, i2] and the relevant byte address of the internal data is given by

```
offset = array.strides[0] * i1 + array.strides[1] * i2
```

- When reshaping an array, NumPy avoids copies when possible by modifying the strides attribute.
- For example, when transposing a matrix, the order of strides is reversed, but the underlying data remains identical.
- However, flattening a transposed array cannot be accomplished simply by modifying strides (try it!), so a copy is needed.

- Internal array layout can also explain some unexpected performance discrepancies between very similar NumPy operations.
- As a small exercise, can you explain the following benchmarks?

```
a = np.random.rand(5000, 5000)
%timeit a[0,:].sum()
%timeit a[:,0].sum()
100000 loops, best of 3: 9.57 μs per loop
10000 loops, best of 3: 68.3 μs per loop
```



Adapted from:

http://materials.jeremybejarano.com/MPIwithPython/introMPI.html https://mpi4py.readthedocs.io/en/stable/intro.html#what-is-mpi

MPI

- MPI, the Message Passing Interface, is a standardized and portable message-passing system designed to function on a wide variety of parallel computers.
- The standard defines the syntax and semantics of library routines and allows users to write portable programs in the main scientific programming languages (Fortran, C, or C++).
- Since its release, the MPI specification has become the leading standard for message-passing libraries for parallel computers.
- Implementations are available from vendors of high-performance computers and from well known open source projects like MPICH, Open MPI or LAM.

Introduction to MPI

- As tradition has it, we will introduce you to MPI programming using a variation on the standard hello world program.
- Our first MPI python program will be the Hello World program for multiple processes. The source code is as follows:

```
#hello.py
from mpi4py import MPI
comm = MPI.COMM_WORLD
rank = comm.Get_rank()
print("hello world from process ", rank)
```

 After saving this text as hello.py, it is executed using the following command-line syntax, run from the file's directory:

```
$ mpiexec -n 5 python hello.py
```

 The above command will execute five python processes which can all communicate with each other.

Introduction to MPI

• When each program runs, it will print hello, and tell you its rank:

```
hello world from process 0
hello world from process 1
hello world from process 3
hello world from process 2
hello world from process 4
```

- Notice that when you try this on your own, they do not necessarily print in order. This is because 5 separate processes are running on different processors, and we cannot know beforehand which one will execute its print statement first.
- If the processes are being scheduled on the same processor instead of multiple processors, then it is up to the operating system to schedule the processes, and it has no preference of any one of our processes over any other process of ours. In essence, each process executes autonomously.

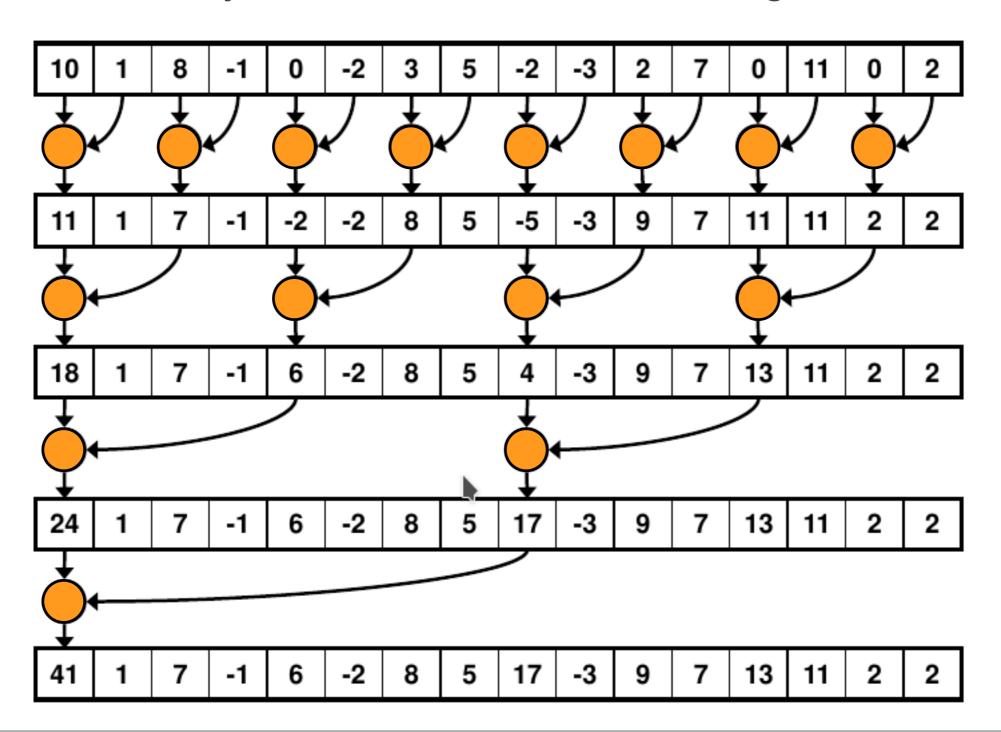
Point To Point Communication

- The simplest message passing involves two processes: a sender and a receiver.
- Let's make two processes. One will draw a random number and then send it to the other.
- We will do this using the routines Comm.Send and Comm.Recv:

```
#pass random draw.py
#example to run: mpiexec -n 2 python pass_random draw.py
import numpy
from mpi4py import MPI
comm = MPI.COMM WORLD
rank = comm.Get rank()
randNum = numpy.zeros(1)
if rank == 1:
        randNum = numpy.random.random sample(1)
        print("Process", rank, "drew the number", randNum[0])
        comm.Send(randNum, dest=0)
if rank == 0:
        print("Process", rank, "before receiving has the number", randNum[0])
        comm.Recv(randNum, source=1)
        print("Process", rank, "received the number", randNum[0])
```

Collective Communication

- Suppose we have eight processes, each with a number to be summed.
- What's the best way to do the sum while minimising communication?



Collective Communication

```
#collective.py
#example to run: mpiexec -n 4 python collective.py 10000
import numpy
import sys
from mpi4py import MPI
comm = MPI.COMM WORLD
rank = comm.Get rank()
size = comm.Get size()
n = int(sys.argv[1])
x = numpy.linspace(start=float(rank)/size, stop=float(1+rank)/size, num=n//size,
endpoint=False)
cosx = numpy.cos(x)
#initializing variables. mpi4py requires that we pass numpy objects.
integral = numpy.zeros(1)
total = numpy.zeros(1)
# perform local computation. Each process integrates its own interval
integral[0] = numpy.sum(cosx)*1.0/n
print("Estimate of integral of cos(x) from %f to %f is %f" % (float(rank)/size,
float(1+rank)/size, integral))
```

Collective Communication

```
# communication
# root node receives results with a collective "reduce"
comm.Reduce(integral, total, op=MPI.SUM, root=0)

# root process prints results
if comm.rank == 0:
    print("With n=%d our estimate of the integral from 0 to 1 of cos(x)
is %f" % (n, total))
    print("Exact integral (sin(1)) is %f" % (numpy.sin(1.0)))
```

More on MPI

- There is a wealth of resources about MPI and the subject is too vast for one lecture
- Just be aware that you ca use it from Python with mpi4py
- For more information about MPI in general check:
 - http://mpitutorial.com/
- And for mpi4py in particular check the documentation at:
 - https://mpi4py.readthedocs.io/en/stable/index.html



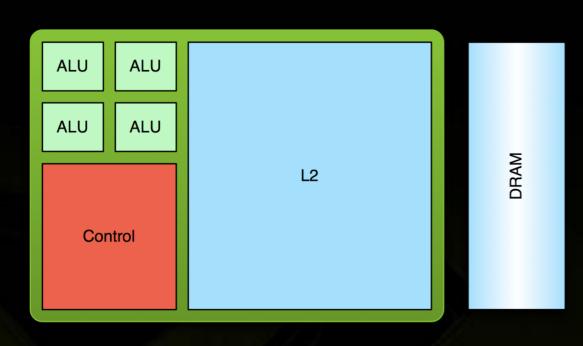
Adapted from:

http://www.cc.gatech.edu/~vetter/keeneland/tutorial-2011-04-14/02-cuda-overview.pdf https://documen.tician.de/pycuda/tutorial.html

- Another way to obtain large gains in performance is to use GPUs to carry out a large fraction of the calculations
- GPUs have enormous computing power. For example an NVIDIA RTX 2090 is capable of 36 TFLOPS in single precision for \$1500.
- For comparison an AMD Epic 7742 (a top of the line \$5000 CPU) reaches only 3.4 TFLOPS
- The two main languages to program GPUs are OpenCL and CUDA
- You can use them from Python using the packages PyCuda and PyOpenCL, but this requires knowledge CUDA and OpenCL respectively.
- Another way to take advantage of them is to use libraries to do their computations in GPUs.
- An example is CuPy which aims to be NumPy on GPUs.

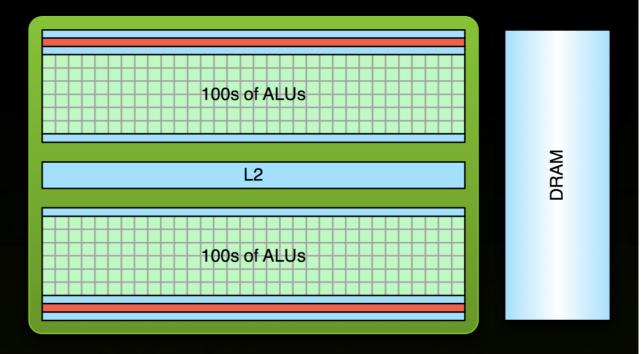
GPU

Low Latency or High Throughput?

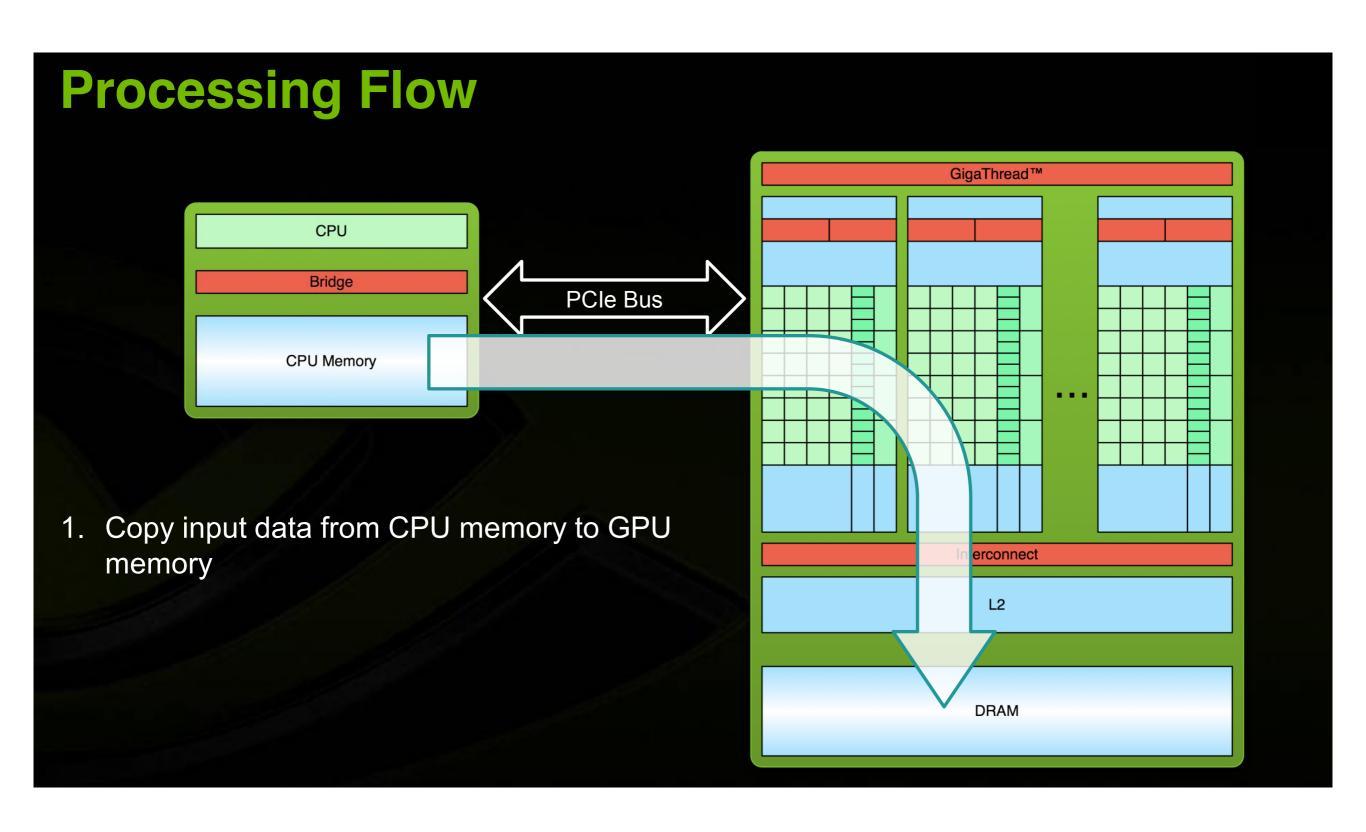


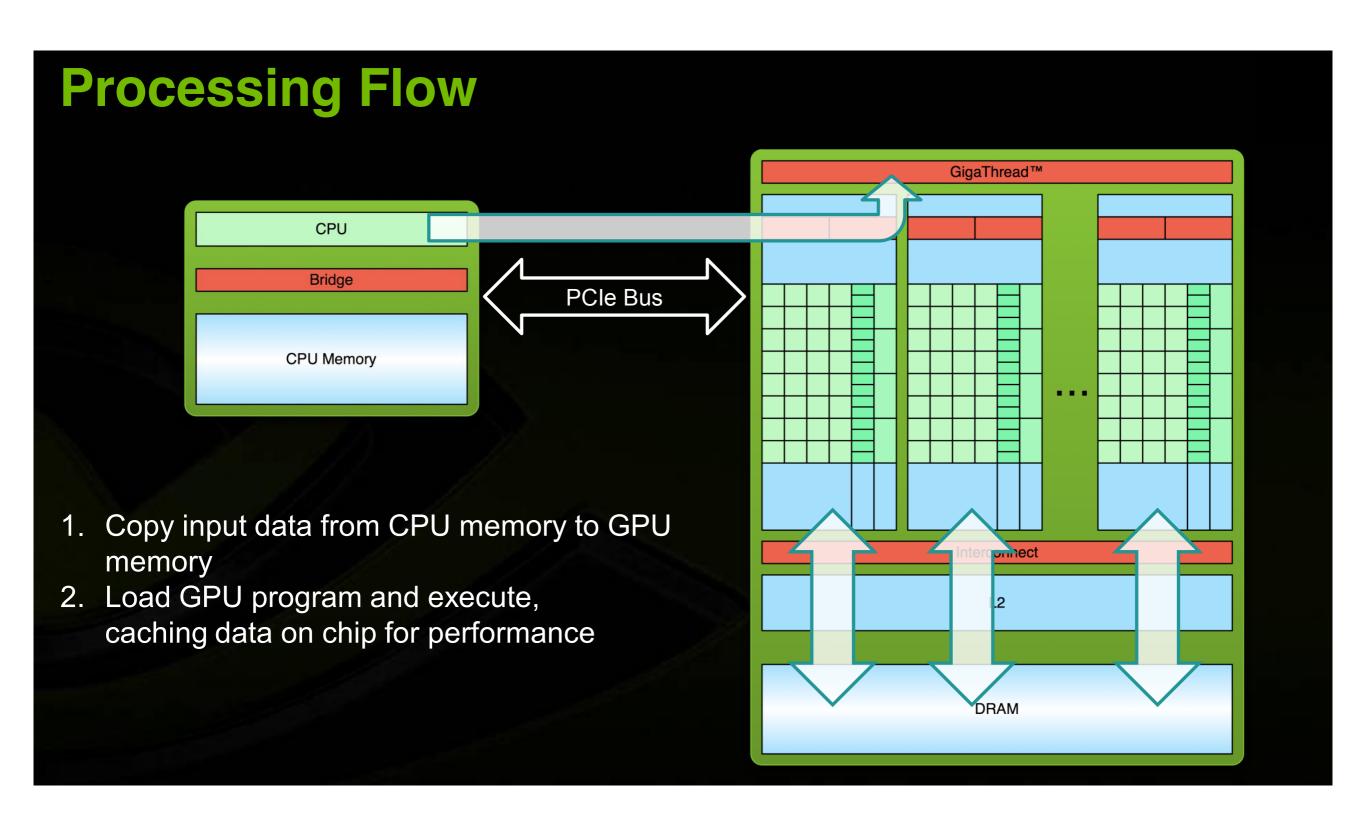
CPU

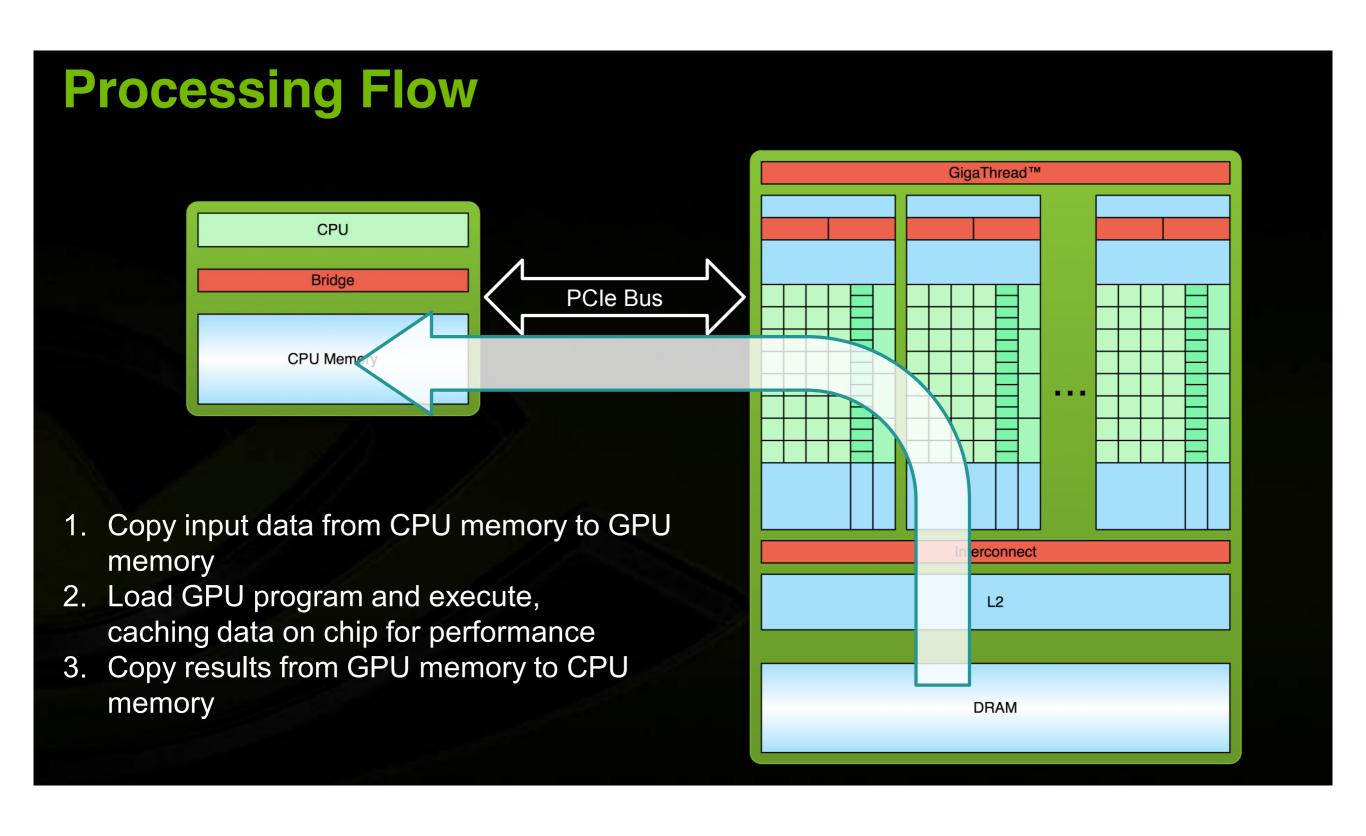
- Optimized for low-latency access to cached data sets
- Control logic for out-of-order and speculative execution



- Optimized for data-parallel, throughput computation
- Architecture tolerant of memory latency
- More transistors dedicated to computation







PyCUDA example

```
import pycuda.driver as cuda
import pycuda.autoinit
from pycuda.compiler import SourceModule
import numpy
# Let's make a 4x4 array of random numbers
a = numpy.random.randn(4,4)
# But wait, a consists of double precision numbers,
# but most nVidia devices only support single precision:
a = a.astype(numpy.float32)
# Finally, we need somewhere to transfer data to,
# so we need to allocate memory on the device
a gpu = cuda.mem alloc(a.nbytes)
# As a last step, we need to transfer the data to the GPU:
cuda.memcpy htod(a gpu, a)
# Kernel to double each entry in an array
# which is compiled by SourceModule
mod = SourceModule("""
    global void doublify(float *a)
    int idx = threadIdx.x + threadIdx.y*4;
    a[idx] *= 2;
  """)
```

PyCUDA example

```
# Grab the compiled function
func = mod.get function("doublify")
# Execute the function on the GPU
# Using one block with 4x4 threads
func(a gpu, block=(4,4,1))
# Finally, we fetch the data back from the GPU and
# display it, together with the original a
a doubled = numpy.empty like(a)
cuda.memcpy_dtoh(a_doubled, a_gpu)
print(a doubled)
print(a)
[[ 0.51360393    1.40589952    2.25009012    3.02563429]
[-0.75841576 -1.18757617 2.72269917 3.12156057]
 0.28826082 -2.92448163
                          1.21624792 2.86353827]
[ 1.57651746  0.63500965  2.21570683  -0.44537592]]
[ 0.25680196 0.70294976 1.12504506 1.51281714]
 [-0.37920788 -0.59378809 1.36134958 1.56078029]
 0.14413041 -1.46224082
                          0.60812396 1.43176913
 [ 0.78825873  0.31750482  1.10785341  -0.22268796]]
```

PyCUDA example

- Now lets do it the easy way!
- PyCUDA has some support for NumPy arrays so you don't need to write kernels for many things!

```
import pycuda.gpuarray as gpuarray
import pycuda.driver as cuda
import pycuda.autoinit
import numpy

a_gpu = gpuarray.to_gpu(numpy.random.randn(4,4).astype(numpy.float32))
a_doubled = (2*a_gpu).get()
print(a_doubled)
print(a_gpu)
```

For more information check the pycuda.gpuarray.GPUArray class

CuPy

- NVIDIA aimed to make GPUs easy to use from Python by supporting the development of CuPy, a NumPy-compatible array library accelerated by CUDA.
- It is Open Source (https://github.com/cupy/cupy), but only supports
 CUDA.

```
# cupy pi.py
                                                # numpy_pi.py
import cupy as np
                                                import numpy as np
# Monte Carlo estimation of pi
                                                # Monte Carlo estimation of pi
def calc pi device(samples):
                                                def calc_pi_device(samples):
   x = np.random.random(size=samples)
                                                    x = np.random.random(size=samples)
   y = np.random.random(size=samples)
                                                    y = np.random.random(size=samples)
                                                    within_unit_circle = (x * x + y * y) < 1.0
   within unit circle = (x * x + y * y) < 1.0
                                                    return 4.0 * sum(within_unit_circle) /
    return 4.0 * np.sum(within unit circle) /
                                                samples
samples
  In [7]: %timeit numpy_pi.calc_pi_device(100000)
  210 ms ± 33.3 ms per loop (mean ± std. dev. of 7 runs, 10 loops each)
  In [8]: %timeit cupy_pi.calc_pi_device(100000)
  242 \mu s \pm 9.38 \mu s per loop (mean \pm std. dev. of 7 runs, 1000 loops each)
```

MPI and **GPUs**

For what kind of tasks are GPUs better suited than CPUs? What about MPI?

References

code examples have been taken and adapted from

- https://github.com/jrjohansson/scientific-python-lectures/blob/master/ Lecture-2-Numpy.ipynb
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- http://scipy-cookbook.readthedocs.io/items/ViewsVsCopies.html
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