

ADVANCED NUMPY

In [193]:

```
import numpy as np
import pandas as pd
import numba as nb
```

In []:

```
# Advanced NumPy

# Let's go deeper into the NumPy library for array computing.
# This will include more internal detail about the ndarray type
# and more advanced array manipulations and algorithms.
```

In []:

```
# Ndarray Object Internals

# The NumPy ndarray provides a means to interpret a block of homogeneous data (either contiguous or strided)
# as a multidimensional array object.
# The data type, or dtype, determines how the data is interpreted as being floating point, integer,
# boolean, or any of the other types.

# Part of what makes ndarray flexible is that every array object is a strided view on a block of data.
# You might wonder, for example, how the array view arr[:,2, :-1] does not copy any data.
# The reason is that the ndarray is more than just a chunk of memory and a dtype;
# it also has “striding” information that enables the array to move through memory with varying step sizes.
# More precisely, the ndarray internally consists of the following:

# • A pointer to data—that is, a block of data in RAM or in a memory-mapped file
# • The data type or dtype, describing fixed-size value cells in the array
# • A tuple indicating the array’s shape
# • A tuple of strides, integers indicating the number of bytes to “step” in order to
# advance one element along a dimension
```

In [3]:

```
# For example, a 10 × 5 array would have shape (10, 5):  
  
np.ones((10, 5)).shape
```

Out[3]:

(10, 5)

In [4]:

```
# A typical (C order) 3 × 4 × 5 array of float64 (8-byte) values has strides (160, 40, 8)  
# (knowing about the strides can be useful because, in general, the larger the strides  
# on a particular axis,  
# the more costly it is to perform computation along that axis):  
  
np.ones((3, 4, 5), dtype=np.float64).strides
```

Out[4]:

(160, 40, 8)

In []:

```
# While it is rare that a typical NumPy user would be interested in the array strides,  
# they are the critical ingredient in constructing “zero-copy” array views. Strides can  
# even be negative,  
# which enables an array to move “backward” through memory  
# (this would be the case, for example, in a slice like obj[::-1] or obj[:, ::-1]).
```

In [6]:

```
# NumPy dtype Hierarchy  
  
# You may occasionally have code that needs to check whether an array contains integers,  
# floating-point numbers, strings, or Python objects.  
# Because there are multiple types of floating-point numbers (float16 through float128),  
# checking that the dtype is among a list of types would be very verbose.  
# Fortunately, the dtypes have superclasses such as np.integer and np.floating,  
# which can be used in conjunction with the np.issubdtype function:  
  
ints = np.ones(10, dtype=np.uint16)  
  
ints
```

Out[6]:

array([1, 1, 1, 1, 1, 1, 1, 1, 1, 1], dtype=uint16)

In [7]:

```
floats = np.ones(10, dtype=np.float32)

floats
```

Out[7]:

```
array([1., 1., 1., 1., 1., 1., 1., 1., 1., 1.], dtype=float32)
```

In [8]:

```
np.issubdtype(ints.dtype, np.integer)
```

Out[8]:

```
True
```

In [9]:

```
np.issubdtype(floats.dtype, np.floating)
```

Out[9]:

```
True
```

In [10]:

```
# You can see all of the parent classes of a specific dtype by calling the type's mro method:

np.float64.mro()
```

Out[10]:

```
[numpy.float64,
 numpy.floating,
 numpy.inexact,
 numpy.number,
 numpy.generic,
 float,
 object]
```

In [11]:

```
# Therefore, we also have:

np.issubdtype(ints.dtype, np.number)
```

Out[11]:

```
True
```

In []:

```
# Most NumPy users will never have to know about this, but it occasionally comes in handy.
```

In []:

```
# Advanced Array Manipulation
```

```
# There are many ways to work with arrays beyond fancy indexing, slicing, and boolean subsetting.
```

```
# While much of the heavy lifting for data analysis applications is handled by higher-level functions
```

```
# in pandas, you may at some point need to write a data algorithm that is not found in one of the existing libraries.
```

In [12]:

```
# Reshaping Arrays
```

```
# In many cases, you can convert an array from one shape to another without copying any data.
```

```
# To do this, pass a tuple indicating the new shape to the reshape array instance method.
```

```
# For example, suppose we had a one-dimensional array of values that we wished to rearrange into a matrix:
```

```
arr = np.arange(8)
```

```
arr
```

Out[12]:

```
array([0, 1, 2, 3, 4, 5, 6, 7])
```

In [17]:

```
arr.reshape((4,2))
```

Out[17]:

```
array([[0, 1],
       [2, 3],
       [4, 5],
       [6, 7]])
```

In [18]:

```
# A multidimensional array can also be reshaped:
```

```
arr.reshape((4, 2)).reshape((2, 4))
```

Out[18]:

```
array([[0, 1, 2, 3],
       [4, 5, 6, 7]])
```

In [19]:

```
# One of the passed shape dimensions can be -1, in which case the value used for the dimension
```

```
# will be inferred from the data:
```

```
arr = np.arange(15)
```

```
arr
```

Out[19]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11, 12, 13, 14])
```

In [20]:

```
arr.reshape((5, -1))
```

Out[20]:

```
array([[ 0,  1,  2],
       [ 3,  4,  5],
       [ 6,  7,  8],
       [ 9, 10, 11],
       [12, 13, 14]])
```

In [22]:

```
# Since an array's shape attribute is a tuple, it can be passed to reshape, too:
```

```
other_arr = np.ones((3, 5))
```

```
other_arr
```

Out[22]:

```
array([[1., 1., 1., 1., 1.],
       [1., 1., 1., 1., 1.],
       [1., 1., 1., 1., 1.]])
```

In [23]:

```
other_arr.shape
```

Out[23]:

```
(3, 5)
```

In [24]:

```
arr.reshape(other_arr.shape)
```

Out[24]:

```
array([[ 0,  1,  2,  3,  4],
       [ 5,  6,  7,  8,  9],
       [10, 11, 12, 13, 14]])
```

In [25]:

```
# The opposite operation of reshape from one-dimensional to a higher dimension  
# is typically known as flattening or raveling:
```

```
arr = np.arange(15).reshape((5, 3))
```

```
arr
```

Out[25]:

```
array([[ 0,  1,  2],
       [ 3,  4,  5],
       [ 6,  7,  8],
       [ 9, 10, 11],
       [12, 13, 14]])
```

In [26]:

```
arr.ravel()
```

Out[26]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11, 12, 13, 14])
```

In [27]:

```
# ravel does not produce a copy of the underlying values if the values in the result  
# were contiguous in the original array.  
# The flatten method behaves like ravel except it always returns a copy of the data:
```

```
arr.flatten()
```

Out[27]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11, 12, 13, 14])
```

In []:

```
# The data can be reshaped or raveled in different orders. This is a slightly nuanced topic  
# for new NumPy users  
# and is therefore the next subtopic.
```

In [28]:

```
# C Versus Fortran Order

# NumPy gives you control and flexibility over the layout of your data in memory.
# By default, NumPy arrays are created in row major order.
# Spatially this means that if you have a two-dimensional array of data,
# the items in each row of the array are stored in adjacent memory locations.
# The alternative to row major ordering is column major order,
# which means that values within each column of data are stored in adjacent memory locations.

# For historical reasons, row and column major order are also known as C and Fortran order, respectively.
# In the FORTRAN 77 language, matrices are all column major.

# Functions like reshape and ravel accept an order argument indicating the order to use the data in the array.
# This is usually set to 'C' or 'F' in most cases
# (there are also less commonly used options 'A' and 'K'; see the NumPy documentation):

arr = np.arange(12).reshape((3, 4))

arr
```

Out[28]:

```
array([[ 0,  1,  2,  3],
       [ 4,  5,  6,  7],
       [ 8,  9, 10, 11]])
```

In [29]:

```
arr.ravel()
```

Out[29]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11])
```

In [30]:

```
arr.ravel('C')
```

Out[30]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11])
```

In [31]:

```
arr.ravel('F')
```

Out[31]:

```
array([ 0,  4,  8,  1,  5,  9,  2,  6, 10,  3,  7, 11])
```

In [32]:

```
arr.ravel('A')
```

Out[32]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11])
```

In [33]:

```
arr.ravel('K')
```

Out[33]:

```
array([ 0,  1,  2,  3,  4,  5,  6,  7,  8,  9, 10, 11])
```

In []:

```
# Reshaping arrays with more than two dimensions can be a bit mind-bending.  
# The key difference between C and Fortran order is the way in which the dimensions are walked:  
  
#           C/row major order  
#           Traverse higher dimensions first (e.g., axis 1 before advancing on axis 0).  
  
#           Fortran/column major order  
#           Traverse higher dimensions last (e.g., axis 0 before advancing on axis 1).
```

In [34]:

```
# Concatenating and Splitting Arrays  
  
# numpy.concatenate takes a sequence (tuple, list, etc.) of arrays and joins them together  
# in order along the input axis:  
  
arr1 = np.array([[1, 2, 3], [4, 5, 6]])  
  
arr1
```

Out[34]:

```
array([[1, 2, 3],  
       [4, 5, 6]])
```

In [36]:

```
arr2 = np.array([[7, 8, 9], [10, 11, 12]])  
  
arr2
```

Out[36]:

```
array([[ 7,  8,  9],  
       [10, 11, 12]])
```


In [37]:

```
np.concatenate([arr1, arr2], axis=0)
```

Out[37]:

```
array([[ 1,  2,  3],
       [ 4,  5,  6],
       [ 7,  8,  9],
       [10, 11, 12]])
```

In [40]:

```
np.concatenate([arr1, arr2], axis=1)
```

Out[40]:

```
array([[ 1,  2,  3,  7,  8,  9],
       [ 4,  5,  6, 10, 11, 12]])
```

In [41]:

```
# There are some convenience functions, like vstack and hstack, for common kinds of concatenation.
```

```
# The preceding operations could have been expressed as:
```

```
np.vstack((arr1, arr2))
```

Out[41]:

```
array([[ 1,  2,  3],
       [ 4,  5,  6],
       [ 7,  8,  9],
       [10, 11, 12]])
```

In [42]:

```
np.hstack((arr1, arr2))
```

Out[42]:

```
array([[ 1,  2,  3,  7,  8,  9],
       [ 4,  5,  6, 10, 11, 12]])
```

In [43]:

```
# split, on the other hand, slices apart an array into multiple arrays along an axis:
```

```
arr = np.random.randn(5, 2)
```

```
arr
```

Out[43]:

```
array([[ 1.33222118, -1.81806039],
       [-0.15346484, -0.13432537],
       [-1.20499111,  1.7143637 ],
       [ 0.29597033, -0.46980145],
       [ 0.28043885,  0.71384464]])
```

In [44]:

```
first, second, third = np.split(arr, [1, 3])
```

```
first
```

Out[44]:

```
array([[ 1.33222118, -1.81806039]])
```

In [45]:

```
second
```

Out[45]:

```
array([[ -0.15346484, -0.13432537],
       [-1.20499111,  1.7143637 ]])
```

In [46]:

```
third
```

Out[46]:

```
array([[ 0.29597033, -0.46980145],
       [ 0.28043885,  0.71384464]])
```

In []:

```
# The value [1, 3] passed to np.split indicate the indices at which to split the array into pieces.
```

```
# Array concatenation functions
```

#	Function	Description
#	<code>concatenate</code>	Most general function, concatenates collection of arrays along one axis
#	<code>vstack, row_stack</code>	Stack arrays row-wise (along axis 0)
#	<code>hstack</code>	Stack arrays column-wise (along axis 1)
#	<code>column_stack</code>	Like <code>hstack</code> , but converts 1D arrays to 2D column vectors first
#	<code>dstack</code>	Stack arrays “depth”-wise (along axis 2)
#	<code>split</code>	Split array at passed locations along a particular axis
#	<code>hsplit/vsplit</code>	Convenience functions for splitting on axis 0 and 1, respectively

In [47]:

```
# Stacking helpers: r_ and c_
```

```
# There are two special objects in the NumPy namespace, r_ and c_, that make stacking arrays more concise:
```

```
arr = np.arange(6)
```

```
arr
```

Out[47]:

```
array([0, 1, 2, 3, 4, 5])
```

In [49]:

```
arr1 = arr.reshape((3, 2))
```

```
arr1
```

Out[49]:

```
array([[0, 1],
       [2, 3],
       [4, 5]])
```

In [50]:

```
arr2 = np.random.randn(3, 2)

arr2
```

Out[50]:

```
array([[ 0.29376076,  0.68752243],
       [-2.53306114,  1.66854045],
       [ 0.01309395,  1.70704312]])
```

In [51]:

```
np.r_[arr1, arr2]
```

Out[51]:

```
array([[ 0.          ,  1.          ],
       [ 2.          ,  3.          ],
       [ 4.          ,  5.          ],
       [ 0.29376076,  0.68752243],
       [-2.53306114,  1.66854045],
       [ 0.01309395,  1.70704312]])
```

In [54]:

```
np.c_[arr1, arr2]
```

Out[54]:

```
array([[ 0.          ,  1.          ,  0.29376076,  0.68752243],
       [ 2.          ,  3.          , -2.53306114,  1.66854045],
       [ 4.          ,  5.          ,  0.01309395,  1.70704312]])
```

In [60]:

```
np.c_[np.r_[arr1, arr2], arr]
```

Out[60]:

```
array([[ 0.          ,  1.          ,  0.          ],
       [ 2.          ,  3.          ,  1.          ],
       [ 4.          ,  5.          ,  2.          ],
       [ 0.29376076,  0.68752243,  3.          ],
       [-2.53306114,  1.66854045,  4.          ],
       [ 0.01309395,  1.70704312,  5.          ]])
```

In [61]:

```
# These additionally can translate slices to arrays:
```

```
np.c_[1:6, -10:-5]
```

Out[61]:

```
array([[ 1, -10],
       [ 2, -9],
       [ 3, -8],
       [ 4, -7],
       [ 5, -6]])
```

In [62]:

```
np.r_[1:6, -10:-5]
```

Out[62]:

```
array([ 1,  2,  3,  4,  5, -10, -9, -8, -7, -6])
```

In []:

```
# See the docstring for more on what you can do with c_ and r_.
```

In [63]:

```
# Repeating Elements: tile and repeat
```

```
# Two useful tools for repeating or replicating arrays to produce larger arrays
```

```
# are the repeat and tile functions.
```

```
# repeat replicates each element in an array some number of times, producing a larger array:
```

```
arr = np.arange(3)
```

```
arr
```

Out[63]:

```
array([0, 1, 2])
```

In [64]:

```
arr.repeat(3)
```

Out[64]:

```
array([0, 0, 0, 1, 1, 1, 2, 2, 2])
```

In [65]:

```
arr.repeat(4)
```

Out[65]:

```
array([0, 0, 0, 0, 1, 1, 1, 1, 2, 2, 2, 2])
```

In [66]:

```
arr.repeat(5)
```

Out[66]:

```
array([0, 0, 0, 0, 0, 1, 1, 1, 1, 1, 2, 2, 2, 2, 2])
```

In [69]:

```
# The need to replicate or repeat arrays can be less common with NumPy than it is with other array programming frameworks like MATLAB. One reason for this is that broadcasting often fills this need better.
```

```
# By default, if you pass an integer, each element will be repeated that number of times.
# If you pass an array of integers, each element can be repeated a different number of times:
```

```
arr.repeat([2, 3, 4])
```

Out[69]:

```
array([0, 0, 1, 1, 1, 2, 2, 2, 2])
```

In [70]:

```
arr.repeat([4, 7, 9])
```

Out[70]:

```
array([0, 0, 0, 0, 1, 1, 1, 1, 1, 1, 1, 1, 2, 2, 2, 2, 2, 2, 2, 2, 2])
```

In [71]:

```
# Multidimensional arrays can have their elements repeated along a particular axis.
```

```
arr = np.random.randn(2, 2)
```

```
arr
```

Out[71]:

```
array([[ -0.64286408, -0.88167946],
       [-0.03991212, -1.66391726]])
```

In [72]:

```
arr.repeat(2, axis=0)
```

Out[72]:

```
array([[ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726]])
```

In [73]:

```
arr.repeat(2, axis=1)
```

Out[73]:

```
array([[ -0.64286408, -0.64286408, -0.88167946, -0.88167946],
       [ -0.03991212, -0.03991212, -1.66391726, -1.66391726]])
```

In [74]:

```
arr.repeat(7, axis=0)
```

Out[74]:

```
array([[ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726],
       [ -0.03991212, -1.66391726]])
```

In [75]:

```
arr.repeat(7, axis=1)
```

Out[75]:

```
array([[ -0.64286408, -0.64286408, -0.64286408, -0.64286408, -0.64286408,
        -0.64286408, -0.64286408, -0.88167946, -0.88167946, -0.88167946,
        -0.88167946, -0.88167946, -0.88167946, -0.88167946],
       [ -0.03991212, -0.03991212, -0.03991212, -0.03991212, -0.03991212,
        -0.03991212, -0.03991212, -1.66391726, -1.66391726, -1.66391726,
        -1.66391726, -1.66391726, -1.66391726, -1.66391726]])
```

In [76]:

```
# Note that if no axis is passed, the array will be flattened first, which is likely not what you want.  
# Similarly, you can pass an array of integers when repeating a multidimensional array  
# to repeat a given slice a different number of times:
```

```
arr.repeat([2, 3], axis=0)
```

Out[76]:

```
array([[ -0.64286408, -0.88167946],  
       [ -0.64286408, -0.88167946],  
       [ -0.03991212, -1.66391726],  
       [ -0.03991212, -1.66391726],  
       [ -0.03991212, -1.66391726]])
```

In [77]:

```
arr.repeat([2, 3], axis=1)
```

Out[77]:

```
array([[ -0.64286408, -0.64286408, -0.88167946, -0.88167946, -0.88167946],  
       [ -0.03991212, -0.03991212, -1.66391726, -1.66391726, -1.66391726]])
```

In [78]:

```
# tile, on the other hand, is a shortcut for stacking copies of an array along an axis.  
# Visually you can think of it as being akin to "laying down tiles":
```

```
arr
```

Out[78]:

```
array([[ -0.64286408, -0.88167946],  
       [ -0.03991212, -1.66391726]])
```

In [79]:

```
np.tile(arr, 2)
```

Out[79]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946],  
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```


In [80]:

```
np.tile(arr, 7)
```

Out[80]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946, -0.64286408,
        -0.88167946, -0.64286408, -0.88167946, -0.64286408, -0.88167946,
        -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [-0.03991212, -1.66391726, -0.03991212, -1.66391726, -0.03991212,
        -1.66391726, -0.03991212, -1.66391726, -0.03991212, -1.66391726,
        -0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```

In [81]:

```
# The second argument is the number of tiles; with a scalar, the tiling is made row by row,
# rather than column by column.
# The second argument to tile can be a tuple indicating the layout of the "tiling":

arr
```

Out[81]:

```
array([[ -0.64286408, -0.88167946],
       [-0.03991212, -1.66391726]])
```

In [82]:

```
np.tile(arr, (2, 1))
```

Out[82]:

```
array([[ -0.64286408, -0.88167946],
       [-0.03991212, -1.66391726],
       [-0.64286408, -0.88167946],
       [-0.03991212, -1.66391726]])
```

In [85]:

```
np.tile(arr, (1, 2))
```

Out[85]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [-0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```

In [86]:

```
np.tile(arr, (3, 2))
```

Out[86]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```

In [87]:

```
np.tile(arr, (2, 3))
```

Out[87]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946, -0.64286408,
        -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726, -0.03991212,
        -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946, -0.64286408,
        -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726, -0.03991212,
        -1.66391726]])
```

In [83]:

```
np.tile(arr, (2, 7))
```

Out[83]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946, -0.64286408,
        -0.88167946, -0.64286408, -0.88167946, -0.64286408, -0.88167946,
        -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726, -0.03991212,
        -1.66391726, -0.03991212, -1.66391726, -0.03991212, -1.66391726,
        -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946, -0.64286408,
        -0.88167946, -0.64286408, -0.88167946, -0.64286408, -0.88167946,
        -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726, -0.03991212,
        -1.66391726, -0.03991212, -1.66391726, -0.03991212, -1.66391726,
        -0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```

In [84]:

```
np.tile(arr, (7, 2))
```

Out[84]:

```
array([[ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726],
       [ -0.64286408, -0.88167946, -0.64286408, -0.88167946],
       [ -0.03991212, -1.66391726, -0.03991212, -1.66391726]])
```

In [88]:

```
# Fancy Indexing Equivalents: take and put
```

```
# One way to get and set subsets of arrays is by fancy indexing using integer arrays:
```

```
arr = np.arange(10) * 100
```

```
arr
```

Out[88]:

```
array([  0, 100, 200, 300, 400, 500, 600, 700, 800, 900])
```

In [89]:

```
inds = [7, 1, 2, 6]
```

```
inds
```

Out[89]:

```
[7, 1, 2, 6]
```

In [90]:

```
arr[inds]
```

Out[90]:

```
array([700, 100, 200, 600])
```

In [91]:

```
# There are alternative ndarray methods that are useful in the special case of only making  
# a selection on a single axis:
```

```
arr.take(inds)
```

Out[91]:

```
array([700, 100, 200, 600])
```

In [92]:

```
arr.put(inds, 42)
```

```
arr
```

Out[92]:

```
array([ 0, 42, 42, 300, 400, 500, 42, 42, 800, 900])
```

In [93]:

```
arr.put(inds, [40, 41, 42, 43])
```

```
arr
```

Out[93]:

```
array([ 0, 41, 42, 300, 400, 500, 43, 40, 800, 900])
```

In [95]:

```
# To use take along other axes, you can pass the axis keyword:
```

```
inds = [2, 0, 2, 1]
```

```
inds
```

Out[95]:

```
[2, 0, 2, 1]
```

In [96]:

```
arr = np.random.randn(2, 4)
```

```
arr
```

Out[96]:

```
array([[ -0.6864408 ,  0.07286014,  0.58338652, -2.01298925],  
       [ 0.34688456,  0.14186142,  0.18422074,  0.77167161]])
```

In [97]:

```
arr.take(inds, axis=1)
```

Out[97]:

```
array([[ 0.58338652, -0.6864408 ,  0.58338652,  0.07286014],
       [ 0.18422074,  0.34688456,  0.18422074,  0.14186142]])
```

In [99]:

```
# put does not accept an axis argument but rather indexes into the flattened (onedimensional, C order) version
# of the array.
# Thus, when you need to set elements using an index array on other axes,
# it is often easiest to use fancy indexing.
```

In [6]:

```
# Broadcasting

# Broadcasting describes how arithmetic works between arrays of different shapes.
# It can be a powerful feature, but one that can cause confusion, even for experienced
users.
# The simplest example of broadcasting occurs when combining a scalar value with an array:

arr = np.arange(5)

arr
```

Out[6]:

```
array([0, 1, 2, 3, 4])
```

In [7]:

```
arr * 4
```

Out[7]:

```
array([ 0,  4,  8, 12, 16])
```

In [8]:

```
# Above we say that the scalar value 4 has been broadcast to all of the other elements in  
# the multiplication operation.
```

```
# For example, we can demean each column of an array by subtracting the column means.  
# In this case, it is very simple:
```

```
arr = np.random.randn(4, 3)
```

```
arr
```

Out[8]:

```
array([[ -0.4290365,  -1.23478276,   0.04131859],  
       [  0.41796202,  -1.44762384,   0.42056144],  
       [  0.00224038,   0.47297833,  -0.26550207],  
       [-0.49401702,  -1.02432231,   0.52866203]])
```

In [9]:

```
arr.mean()
```

Out[9]:

```
-0.2509634766353159
```

In [10]:

```
arr.mean(0)
```

Out[10]:

```
array([-0.12571278, -0.80843765,   0.18126   ])
```

In [11]:

```
arr.mean(1)
```

Out[11]:

```
array([-0.54083356, -0.20303346,   0.06990555, -0.32989243])
```

In [13]:

```
demeaned = arr - arr.mean(0)
```

```
demeaned
```

Out[13]:

```
array([[ -0.30332372,  -0.42634511,  -0.13994141],  
       [  0.5436748 ,  -0.6391862 ,   0.23930144],  
       [  0.12795316,   1.28141597,  -0.44676207],  
       [-0.36830424,  -0.21588466,   0.34740203]])
```

In [15]:

```
demeaned.mean()
```

Out[15]:

```
-1.3877787807814457e-17
```

In [16]:

```
demeaned.mean(0)
```

Out[16]:

```
array([ 1.38777878e-17, -2.77555756e-17,  0.00000000e+00])
```

In [17]:

```
demeaned.mean(1)
```

Out[17]:

```
array([-0.28987008,  0.04793002,  0.32086902, -0.07892896])
```

In []:

```
# Demeaning the rows as a broadcast operation requires a bit more care.
# Fortunately, broadcasting potentially lower dimensional values across any dimension
of an array
# (like subtracting the row means from each column of a two-dimensional array)
# is possible as long as you follow the rules.

# This brings us to:

#
# The Broadcasting Rule
# Two arrays are compatible for broadcasting if for each trailing dimension (i.e., starting
from the end)
# the axis lengths match or if either of the lengths is 1. Broadcasting is then performed
over the missing
# or length 1 dimensions.
```

In [18]:

```
# Even as an experienced NumPy user, I often find myself having to pause and draw a diagram as I think about  
# the broadcasting rule.  
# Consider the last example and suppose we wished instead to subtract the mean value from each row.  
# Since arr.mean(0) has length 3, it is compatible for broadcasting across axis 0 because the trailing  
# dimension in arr is 3 and therefore matches.  
# According to the rules, to subtract over axis 1 (i.e., subtract the row mean from each row),  
# the smaller array must have shape (4, 1):  
  
arr
```

Out[18]:

```
array([[ -0.4290365,  -1.23478276,   0.04131859],  
       [  0.41796202,  -1.44762384,   0.42056144],  
       [  0.00224038,   0.47297833,  -0.26550207],  
       [-0.49401702,  -1.02432231,   0.52866203]])
```

In [21]:

```
row_means = arr.mean(1)  
  
row_means
```

Out[21]:

```
array([-0.54083356, -0.20303346,  0.06990555, -0.32989243])
```

In [22]:

```
row_means.shape
```

Out[22]:

```
(4,)
```

In [23]:

```
row_means.reshape((4,1))
```

Out[23]:

```
array([[ -0.54083356],  
       [-0.20303346],  
       [ 0.06990555],  
       [-0.32989243]])
```


In [25]:

```
demeaned = arr - row_means.reshape((4,1))  
  
demeaned
```

Out[25]:

```
array([[ 0.11179706, -0.6939492,  0.58215215],  
       [ 0.62099549, -1.24459038,  0.6235949 ],  
       [-0.06766516,  0.40307278, -0.33540762],  
       [-0.16412459, -0.69442987,  0.85855446]])
```

In [26]:

```
demeaned.mean()
```

Out[26]:

```
1.850371707708594e-17
```

In [27]:

```
demeaned.mean(0)
```

Out[27]:

```
array([ 0.1252507, -0.55747417,  0.43222347])
```

In [28]:

```
demeaned.mean(1)
```

Out[28]:

```
array([-3.70074342e-17,  3.70074342e-17,  0.00000000e+00,  0.00000000e+00])
```

In [29]:

```
# Broadcasting Over Other Axes  
  
# Broadcasting with higher dimensional arrays can seem even more mind-bending,  
# but it is really a matter of following the rules. If you don't, you'll get an error  
like this:  
  
arr - arr.mean(1)
```

```
-----  
ValueError                                Traceback (most recent call last)  
<ipython-input-29-9e4a803024ca> in <module>  
      4      # butvit is really a matter of following the rules. If you don't,  
      you'll get an error like this:  
      5  
----> 6 arr - arr.mean(1)
```

ValueError: operands could not be broadcast together with shapes (4,3) (4,)

In [30]:

```
# It's quite common to want to perform an arithmetic operation with a lower dimensional array  
# across axes other than axis 0.  
# According to the broadcasting rule, the "broadcast dimensions" must be 1 in the smaller array.  
# In the example of row demeaning shown here, this meant reshaping the row means to be shape (4, 1)  
# instead of (4,):  
  
arr - arr.mean(1).reshape((4, 1))
```

Out[30]:

```
array([[ 0.11179706, -0.6939492,  0.58215215],  
       [ 0.62099549, -1.24459038,  0.6235949 ],  
       [-0.06766516,  0.40307278, -0.33540762],  
       [-0.16412459, -0.69442987,  0.85855446]])
```

In [32]:

```
# In the three-dimensional case, broadcasting over any of the three dimensions is only a matter of reshaping  
# the data to be shape-compatible.  
  
# A common problem, therefore, is needing to add a new axis with length 1 specifically  
# for broadcasting purposes.  
# Using reshape is one option, but inserting an axis requires constructing a tuple  
# indicating the new shape. This can often be a tedious exercise.  
# Thus, NumPy arrays offer a special syntax for inserting new axes by indexing.  
# We use the special np.newaxis attribute along with "full" slices to insert the new axis:  
  
arr = np.zeros((4, 4))  
  
arr
```

Out[32]:

```
array([[0., 0., 0., 0.],  
       [0., 0., 0., 0.],  
       [0., 0., 0., 0.],  
       [0., 0., 0., 0.]])
```

In [34]:

```
arr_3d = arr[:, np.newaxis, :]  
arr_3d
```

Out[34]:

```
array([[[0., 0., 0., 0.]],  
       [[0., 0., 0., 0.]],  
       [[0., 0., 0., 0.]],  
       [[0., 0., 0., 0.]])
```

In [35]:

```
arr_3d.shape
```

Out[35]:

```
(4, 1, 4)
```

In [37]:

```
arr_1d = np.random.normal(size=3)  
arr_1d
```

Out[37]:

```
array([-0.40719218, -1.32142006,  0.12264432])
```

In [38]:

```
arr_1d[:, np.newaxis]
```

Out[38]:

```
array([[-0.40719218],  
       [-1.32142006],  
       [ 0.12264432]])
```

In [39]:

```
arr_1d[np.newaxis, :]
```

Out[39]:

```
array([[-0.40719218, -1.32142006,  0.12264432]])
```

In [40]:

```
# Thus, if we had a three-dimensional array and wanted to demean axis 2, say, we would need to write:
```

```
arr = np.random.randn(3, 4, 5)
```

```
arr
```

Out[40]:

```
array([[[ 0.22067351,  3.01315715,  1.3245668 ,  0.73755868,
          0.20639973],
        [-0.6198023 ,  0.66000681, -0.82840268, -1.11622781,
          0.86700759],
        [-0.47165603, -0.71247087, -1.07217153, -0.11514507,
          0.52126238],
        [-1.22809747,  0.45778851,  0.69904621,  1.6742796 ,
          1.07299997]],

       [[ 0.07624852,  0.46179254, -0.82756269, -1.23275627,
          2.32956618],
        [-0.82614275, -0.77474342, -1.77655293, -0.91926351,
          0.50336781],
        [-1.08196146,  0.29968226,  0.96002316, -0.25009993,
         -0.7572684 ],
        [ 0.74146467, -1.00601945,  0.34643083,  1.12784706,
          0.44480431]],

       [[-0.59561434,  0.51490082,  1.87081664, -0.84672811,
          1.33119253],
        [ 0.55605734, -1.23795331, -1.47202384,  0.71790392,
          0.29493959],
        [ 0.55308327, -0.71064273,  0.31722041, -1.27594141,
         -0.42941873],
        [ 0.65510907,  1.57775787,  1.11032321,  0.72963188,
         -0.73807179]])
```

In [41]:

```
depth_means = arr.mean(2)
```

```
depth_means
```

Out[41]:

```
array([[ 1.10047117, -0.20748368, -0.37003622,  0.53520336],
       [ 0.16145766, -0.75866696, -0.16592487,  0.33090548],
       [ 0.45491351, -0.22821526, -0.30913984,  0.66695005]])
```

In [42]:

```
depth_means.shape
```

Out[42]:

```
(3, 4)
```

In [44]:

```
demeaned = arr - depth_means[:, :, np.newaxis]
demeaned
```

Out[44]:

```
array([[[ -0.87979767,  1.91268598,  0.22409563, -0.3629125 ,
          -0.89407144],
        [ -0.41231862,  0.86749049, -0.620919  , -0.90874413,
          1.07449127],
        [ -0.1016198 , -0.34243465, -0.7021353 ,  0.25489115,
          0.8912986  ],
        [ -1.76330084, -0.07741486,  0.16384285,  1.13907624,
          0.5377966  ]],

       [[ -0.08520913,  0.30033488, -0.98902035, -1.39421393,
          2.16810853],
        [ -0.06747579, -0.01607646, -1.01788597, -0.16059655,
          1.26203477],
        [ -0.91603658,  0.46560713,  1.12594804, -0.08417506,
          -0.59134353],
        [  0.41055918, -1.33692494,  0.01552535,  0.79694157,
          0.11389883 ]],

       [[ -1.05052785,  0.05998731,  1.41590313, -1.30164161,
          0.87627902],
        [  0.7842726 , -1.00973805, -1.24380858,  0.94611918,
          0.52315485],
        [  0.86222311, -0.40150289,  0.62636025, -0.96680157,
          -0.12027889],
        [ -0.01184098,  0.91080782,  0.44337316,  0.06268183,
          -1.40502184 ]]])
```

In [45]:

```
demeaned.mean(2)
```

Out[45]:

```
array([[ -4.44089210e-17,  0.00000000e+00, -4.44089210e-17,
          -4.44089210e-17],
       [  8.88178420e-17,  0.00000000e+00,  4.44089210e-17,
          1.11022302e-17],
       [  0.00000000e+00,  0.00000000e+00,  6.66133815e-17,
          0.00000000e+00]])
```

In []:

```
# You might be wondering if there's a way to generalize demeaning over an axis without sacrificing performance.
```

```
# There is, but it requires some indexing gymnastics:
```

```
def demean_axis(arr, axis=0):  
    means = arr.mean(axis)  
  
    # This generalizes things like[:, :, np.newaxis] to N dimensions  
    indexer = [slice(None)] * arr.ndim  
    indexer[axis] = np.newaxis  
    return arr - means[indexer]
```

In [47]:

```
# Setting Array Values by Broadcasting
```

```
# The same broadcasting rule governing arithmetic operations also applies to setting values via array indexing.
```

```
# In a simple case, we can do things like:
```

```
arr = np.zeros((4, 3))
```

```
arr
```

Out[47]:

```
array([[0., 0., 0.],  
       [0., 0., 0.],  
       [0., 0., 0.],  
       [0., 0., 0.]])
```

In [49]:

```
arr[:] = 5
```

```
arr
```

Out[49]:

```
array([[5., 5., 5.],  
       [5., 5., 5.],  
       [5., 5., 5.],  
       [5., 5., 5.]])
```

In [50]:

```
# However, if we had a one-dimensional array of values we wanted to set into the columns of the array,  
# we can do that as long as the shape is compatible:
```

```
col = np.array([1.28, -0.42, 0.44, 1.6])
```

```
col
```

Out[50]:

```
array([ 1.28, -0.42,  0.44,  1.6 ])
```

In [51]:

```
arr[:,] = col[:, np.newaxis]
```

```
arr
```

Out[51]:

```
array([[ 1.28,  1.28,  1.28],  
       [-0.42, -0.42, -0.42],  
       [ 0.44,  0.44,  0.44],  
       [ 1.6 ,  1.6 ,  1.6 ]])
```

In [52]:

```
arr[:2] = [[-1.37], [0.509]]
```

```
arr
```

Out[52]:

```
array([[ -1.37 , -1.37 , -1.37 ],  
       [  0.509,  0.509,  0.509],  
       [  0.44 ,  0.44 ,  0.44 ],  
       [  1.6  ,  1.6  ,  1.6  ]])
```

In []:

```
#Advanced ufunc Usage

# While many NumPy users will only make use of the fast element-wise operations provided
# by the universal functions, there are a number of additional features that occasionally
# can help you
# write more concise code without loops.

# ufunc Instance Methods

# Each of NumPy's binary ufuncs has special methods for performing certain kinds
# of special vectorized operations:

#           Method           Description
#
#           reduce(x)         Aggregate values by successive applications of the operation
#
#           accumulate(x)      Aggregate values, preserving all partial aggregates
#
#           reduceat(x, bins)  "Local" reduce or "group by"; reduce contiguous slices
#           of data            to produce aggregated array
#
#           outer(x, y)        Apply operation to all pairs of elements in x and y;
#                               the resulting array has shape x.shape + y.shape
```

In [53]:

```
# reduce takes a single array and aggregates its values, optionally along an axis,
# by performing a sequence of binary operations.
# For example, an alternative way to sum elements in an array is to use np.add.reduce:

arr = np.arange(10)

arr
```

Out[53]:

```
array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
```

In [54]:

```
np.add.reduce(arr)
```

Out[54]:

45

In [55]:

```
arr.sum()
```

Out[55]:

45

In [57]:

```
# The starting value (0 for add) depends on the ufunc.  
# If an axis is passed, the reduction is performed along that axis.  
# This allows you to answer certain kinds of questions in a concise way.  
# As a less trivial example, we can use np.logical_and to check whether the values  
# in each row of an array are sorted:
```

```
np.random.seed(12346) # for reproducibility
```

```
arr = np.random.randn(5, 5)
```

```
arr
```

Out[57]:

```
array([[ -8.99822478e-02,  7.59372617e-01,  7.48336101e-01,  
        -9.81497953e-01,  3.65775545e-01],  
       [-3.15442628e-01, -8.66135605e-01,  2.78568155e-02,  
        -4.55597723e-01, -1.60189223e+00],  
       [ 2.48256116e-01, -3.21536673e-01, -8.48730755e-01,  
         4.60468309e-04, -5.46459347e-01],  
       [ 2.53915229e-01,  1.93684246e+00, -7.99504902e-01,  
        -5.69159281e-01,  4.89244731e-02],  
       [-6.49092950e-01, -4.79535727e-01, -9.53521432e-01,  
         1.42253882e+00,  1.75403128e-01]])
```

In [64]:

```
arr[:, :2].sort(1) # sort a few rows
```

```
arr[:, :-1] < arr[:, 1:]
```

Out[64]:

```
array([[ True,  True,  True,  True],  
       [False,  True, False, False],  
       [ True,  True,  True,  True],  
       [ True, False,  True,  True],  
       [ True,  True,  True,  True]])
```

In [65]:

```
np.logical_and.reduce(arr[:, :-1] < arr[:, 1:], axis=1)
```

Out[65]:

```
array([ True, False,  True, False,  True])
```

In [66]:

```
# Note that logical_and.reduce is equivalent to the all method.

# accumulate is related to reduce like cumsum is related to sum.
# It produces an array of the same size with the intermediate "accumulated" values:

arr = np.arange(15).reshape((3, 5))

arr
```

Out[66]:

```
array([[ 0,  1,  2,  3,  4],
       [ 5,  6,  7,  8,  9],
       [10, 11, 12, 13, 14]])
```

In [67]:

```
np.add.accumulate(arr, axis=1)
```

Out[67]:

```
array([[ 0,  1,  3,  6, 10],
       [ 5, 11, 18, 26, 35],
       [10, 21, 33, 46, 60]], dtype=int32)
```

In [68]:

```
np.add.accumulate(arr, axis=0)
```

Out[68]:

```
array([[ 0,  1,  2,  3,  4],
       [ 5,  7,  9, 11, 13],
       [15, 18, 21, 24, 27]], dtype=int32)
```

In [69]:

```
np.add.accumulate(arr)
```

Out[69]:

```
array([[ 0,  1,  2,  3,  4],
       [ 5,  7,  9, 11, 13],
       [15, 18, 21, 24, 27]], dtype=int32)
```

In [70]:

```
# outer performs a pairwise cross-product between two arrays:
```

```
arr = np.arange(3).repeat([1, 2, 2])
```

```
arr
```

Out[70]:

```
array([0, 1, 1, 2, 2])
```

In [71]:

```
np.multiply.outer(arr, np.arange(5))
```

Out[71]:

```
array([[0, 0, 0, 0, 0],
       [0, 1, 2, 3, 4],
       [0, 1, 2, 3, 4],
       [0, 2, 4, 6, 8],
       [0, 2, 4, 6, 8]])
```

In [74]:

```
#The output of outer will have a dimension that is the sum of the dimensions of the input s:
```

```
x, y = np.random.randn(3, 4), np.random.randn(5)
```

```
x
```

Out[74]:

```
array([[ 0.73289771,  0.50470465, -0.7892592 ,  0.5391877 ],
       [ 1.29070685,  0.86761856,  0.41133011,  0.44593599],
       [-0.3171888 , -1.04929141,  1.34589315,  0.35600969]])
```

In [75]:

```
y
```

Out[75]:

```
array([-0.09152874, -0.53496417, -0.03601325, -0.25911386, -0.19944861])
```

In [76]:

```
x, y
```

Out[76]:

```
(array([[ 0.73289771,  0.50470465, -0.7892592 ,  0.5391877 ],
        [ 1.29070685,  0.86761856,  0.41133011,  0.44593599],
        [-0.3171888 , -1.04929141,  1.34589315,  0.35600969]]),
 array([-0.09152874, -0.53496417, -0.03601325, -0.25911386, -0.19944861]))
```

In [78]:

```
result = np.subtract.outer(x, y)
```

```
result
```

Out[78]:

```
array([[[[ 0.82442645,  1.26786188,  0.76891096,  0.99201157,
           0.93234632],
         [ 0.59623339,  1.03966882,  0.5407179 ,  0.76381851,
           0.70415326],
         [-0.69773046, -0.25429503, -0.75324595, -0.53014534,
          -0.58981059],
         [ 0.63071644,  1.07415187,  0.57520095,  0.79830156,
           0.7386363 ]],

        [[ 1.38223559,  1.82567102,  1.3267201 ,  1.5498207 ,
           1.49015545],
         [ 0.95914729,  1.40258272,  0.9036318 ,  1.12673241,
           1.06706716],
         [ 0.50285885,  0.94629427,  0.44734335,  0.67044396,
           0.61077871],
         [ 0.53746473,  0.98090016,  0.48194924,  0.70504985,
           0.6453846 ]],

        [[-0.22566006,  0.21777537, -0.28117555, -0.05807494,
          -0.11774019],
         [-0.95776267, -0.51432725, -1.01327817, -0.79017756,
          -0.84984281],
         [ 1.43742189,  1.88085732,  1.3819064 ,  1.60500701,
           1.54534175],
         [ 0.44753843,  0.89097386,  0.39202294,  0.61512355,
           0.5554583 ]]])
```

In [79]:

```
result.shape
```

Out[79]:

```
(3, 4, 5)
```

In [80]:

```
# The Last method, reduceat, performs a “local reduce,” in essence an array groupby operation  
# in which slices of the array are aggregated together.  
# It accepts a sequence of “bin edges” that indicate how to split and aggregate the values:  
  
arr = np.arange(10)  
  
arr
```

Out[80]:

```
array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
```

In [81]:

```
np.add.reduceat(arr, [0, 5, 8])
```

Out[81]:

```
array([10, 18, 17], dtype=int32)
```

In [82]:

```
# The results are the reductions (here, sums) performed over arr[0:5], arr[5:8], and arr[8:].  
# As with the other methods, you can pass an axis argument:  
  
arr = np.multiply.outer(np.arange(4), np.arange(5))  
  
arr
```

Out[82]:

```
array([[ 0,  0,  0,  0,  0],  
       [ 0,  1,  2,  3,  4],  
       [ 0,  2,  4,  6,  8],  
       [ 0,  3,  6,  9, 12]])
```

In [90]:

```
np.add.reduceat(arr, [0, 2, 4], axis=1)
```

Out[90]:

```
array([[ 0,  0,  0],  
       [ 1,  5,  4],  
       [ 2, 10,  8],  
       [ 3, 15, 12]], dtype=int32)
```

In [95]:

```
# Writing New ufuncs in Python

# There are a number of facilities for creating your own NumPy ufuncs.
# The most general is to use the NumPy C API, but let's look at pure Python ufuncs.

# numpy.frompyfunc accepts a Python function along with a specification for the number of
inputs and outputs.
# For example, a simple function that adds element-wise would be specified as:

def add_elements(x, y):
    return x + y
```

In [97]:

```
add_them = np.frompyfunc(add_elements, 2, 1)

add_them
```

Out[97]:

```
<ufunc '?' (vectorized)'>
```

In [98]:

```
add_them(np.arange(8), np.arange(8))
```

Out[98]:

```
array([0, 2, 4, 6, 8, 10, 12, 14], dtype=object)
```

In [100]:

```
# Functions created using frompyfunc always return arrays of Python objects,
# which can be inconvenient.
# Fortunately, there is an alternative (but slightly less featureful) function, numpy.
vectorize,
# that allows you to specify the output type:

add_them = np.vectorize(add_elements, otypes=[np.float64])

add_them
```

Out[100]:

```
<numpy.vectorize at 0xa010275188>
```

In [101]:

```
add_them(np.arange(8), np.arange(8))
```

Out[101]:

```
array([ 0.,  2.,  4.,  6.,  8., 10., 12., 14.])
```

In [109]:

```
# These functions provide a way to create ufunc-like functions,  
# but they are very slow because they require a Python function call to compute each element,  
# which is a lot slower than NumPy's C-based ufunc loops:  
  
arr = np.random.randn(10000)  
  
arr
```

Out[109]:

```
array([ 0.10533048, -0.81101068, -1.14124808, ...,  1.16323222,  
       -0.80433614,  0.08255916])
```

In [112]:

```
%timeit add_them(arr, arr)
```

5.67 ms \pm 723 μ s per loop (mean \pm std. dev. of 7 runs, 100 loops each)

In [113]:

```
%timeit np.add(arr, arr)
```

17.9 μ s \pm 1.05 μ s per loop (mean \pm std. dev. of 7 runs, 100000 loops each)

In [115]:

```
# Structured and Record Arrays  
  
# You may have noticed up until now that ndarray is a homogeneous data container;  
# that is, it represents a block of memory in which each element takes up the same number of bytes,  
# determined by the dtype.  
# On the surface, this would appear to not allow you to represent heterogeneous or tabular-like data.  
# A structured array is an ndarray in which each element can be thought of as representing a struct in C  
# (hence the "structured" name) or a row in a SQL table with multiple named fields:  
  
dtype = [('x', np.float64), ('y', np.int32)]  
  
dtype
```

Out[115]:

```
[('x', numpy.float64), ('y', numpy.int32)]
```

In [116]:

```
sarr = np.array([(1.5, 6), (np.pi, -2)], dtype=dtype)

sarr
```

Out[116]:

```
array([(1.5, 6), (3.14159265, -2)],
      dtype=[('x', '<f8'), ('y', '<i4')])
```

In [117]:

```
# There are several ways to specify a structured dtype (see the online NumPy documentatio
n).
# One typical way is as a list of tuples with (field_name, field_data_type).
# Now, the elements of the array are tuple-like objects whose elements can be accessed
like a dictionary:

sarr[0]
```

Out[117]:

```
(1.5, 6)
```

In [118]:

```
sarr[0]['y']
```

Out[118]:

```
6
```

In [119]:

```
# The field names are stored in the dtype.names attribute.
# When you access a field on the structured array, a strided view on the data is retur
ned,
# thus copying nothing:

sarr['x']
```

Out[119]:

```
array([1.5, 3.14159265])
```


In [120]:

```
# Nested dtypes and Multidimensional Fields

# When specifying a structured dtype, you can additionally pass a shape (as an int or tuple):

dtype = [('x', np.int64, 3), ('y', np.int32)]

dtype
```

Out[120]:

```
[('x', numpy.int64, 3), ('y', numpy.int32)]
```

In [123]:

```
arr = np.zeros(4, dtype=dtype)

arr
```

Out[123]:

```
array([([0, 0, 0], 0), ([0, 0, 0], 0), ([0, 0, 0], 0), ([0, 0, 0], 0)],
      dtype=[('x', '<i8', (3,)), ('y', '<i4')])
```

In [124]:

```
# In this case, the x field now refers to an array of length 3 for each record:

arr[0]['x']
```

Out[124]:

```
array([0, 0, 0], dtype=int64)
```

In [128]:

```
# Conveniently, accessing arr['x'] then returns a two-dimensional array
# instead of a one-dimensional array as in prior examples:

arr['x']
```

Out[128]:

```
array([[0, 0, 0],
       [0, 0, 0],
       [0, 0, 0],
       [0, 0, 0]], dtype=int64)
```

In [130]:

```
# This enables you to express more complicated, nested structures as a single block of memory in an array.  
# You can also nest dtypes to make more complex structures.  
# Here is an example:
```

```
dtype = [('x', [('a', 'f8'), ('b', 'f4')]), ('y', np.int32)]
```

```
dtype
```

Out[130]:

```
[('x', [('a', 'f8'), ('b', 'f4')]), ('y', numpy.int32)]
```

In [131]:

```
data = np.array([(1, 2), 5], [(3, 4), 6]), dtype=dtype)
```

```
data['x']
```

Out[131]:

```
array([(1., 2.), (3., 4.)], dtype=[('a', '<f8'), ('b', '<f4')])
```

In [132]:

```
data['y']
```

Out[132]:

```
array([5, 6])
```

In [133]:

```
data['x']['a']
```

Out[133]:

```
array([1., 3.])
```

In []:

```
# pandas DataFrame does not support this feature directly, though it is similar to hierarchical indexing.
```

In []:

```
# Why Use Structured Arrays?

# Compared with, say, a pandas DataFrame, NumPy structured arrays are a comparatively low-level tool.
# They provide a means to interpreting a block of memory as a tabular structure
# with arbitrarily complex nested columns.
# Since each element in the array is represented in memory as a fixed number of bytes,
# structured arrays provide a very fast and efficient way of writing data to
# and from disk (including memory maps), transporting it over the network, and other such uses.

# As another common use for structured arrays, writing data files as fixed-length
# record byte streams is a common way to serialize data in C and C++ code,
# which is commonly found in legacy systems in industry.
# As long as the format of the file is known
# (the size of each record and the order, byte size, and data type of each element),
# the data can be read into memory with np.fromfile.
```

In [135]:

```
# More About Sorting

# Like Python's built-in list, the ndarray sort instance method is an in-place sort,
# meaning that the array contents are rearranged without producing a new array:

arr = np.random.randn(6)

arr
```

Out[135]:

```
array([-1.4199867 , -0.04596205,  1.27691904,  1.86071261, -0.31313357,
        -0.22521732])
```

In [142]:

```
# When sorting arrays in-place, remember that if the array is a view on a different ndarray,
# the original array will be modified:

arr = np.random.randn(3, 5)

arr
```

Out[142]:

```
array([[ -0.28487369, -0.41144101, -0.39848164, -0.85722451,  0.04447328],
       [-1.92493444, -0.18016988, -0.61207016, -0.02848961, -0.00939751],
       [ 1.16893736,  0.34476488,  0.8829256 , -1.63350964, -0.05559684]])
```

In [146]:

```
arr[:, 0].sort() # Sort first column values in-place  
  
arr
```

Out[146]:

```
array([[ -1.92493444, -0.41144101, -0.39848164, -0.85722451,  0.04447328],  
       [ -0.28487369, -0.18016988, -0.61207016, -0.02848961, -0.00939751],  
       [  1.16893736,  0.34476488,  0.8829256 , -1.63350964, -0.05559684]])
```

In [147]:

```
# On the other hand, numpy.sort creates a new, sorted copy of an array.  
# Otherwise, it accepts the same arguments (such as kind) as ndarray.sort:  
  
arr = np.random.randn(5)  
  
arr
```

Out[147]:

```
array([ 0.23207772,  0.94006741, -0.1233083 ,  0.43651008, -1.0445751 ])
```

In [148]:

```
np.sort(arr)
```

Out[148]:

```
array([-1.0445751 , -0.1233083 ,  0.23207772,  0.43651008,  0.94006741])
```

In [149]:

```
arr
```

Out[149]:

```
array([ 0.23207772,  0.94006741, -0.1233083 ,  0.43651008, -1.0445751 ])
```

In [150]:

```
# All of these sort methods take an axis argument for sorting the sections of data  
# along the passed axis independently:  
  
arr = np.random.randn(3, 5)  
  
arr
```

Out[150]:

```
array([[ 0.37039574, -0.09319317,  0.15047656,  0.60120384,  0.2195671 ],  
       [-2.01196694, -0.69332216, -0.24604222, -1.00853706,  0.52522324],  
       [-2.70506475,  1.24593288,  0.75426158,  0.77936106,  0.96449794]])
```

In [152]:

```
arr.sort(axis=1)

arr
```

Out[152]:

```
array([[ -0.09319317,  0.15047656,  0.2195671 ,  0.37039574,  0.60120384],
       [-2.01196694, -1.00853706, -0.69332216, -0.24604222,  0.52522324],
       [-2.70506475,  0.75426158,  0.77936106,  0.96449794,  1.24593288]])
```

In [153]:

```
arr.sort(axis=0)

arr
```

Out[153]:

```
array([[ -2.70506475, -1.00853706, -0.69332216, -0.24604222,  0.52522324],
       [-2.01196694,  0.15047656,  0.2195671 ,  0.37039574,  0.60120384],
       [-0.09319317,  0.75426158,  0.77936106,  0.96449794,  1.24593288]])
```

In [154]:

```
# You may notice that none of the sort methods have an option to sort in descending order.
# This is a problem in practice because array slicing produces views,
# thus not producing a copy or requiring any computational work.
# Many Python users are familiar with the “trick” that for a list values,
# values[::-1] returns a list in reverse order.
# The same is true for ndarrays:

arr[:, ::-1]
```

Out[154]:

```
array([[ 0.52522324, -0.24604222, -0.69332216, -1.00853706, -2.70506475],
       [ 0.60120384,  0.37039574,  0.2195671 ,  0.15047656, -2.01196694],
       [ 1.24593288,  0.96449794,  0.77936106,  0.75426158, -0.09319317]])
```

In [156]:

```
# Indirect Sorts: argsort and lexsort

# In data analysis you may need to reorder datasets by one or more keys.
# For example, a table of data about some students might need to be sorted by last name,
# then by first name.
# This is an example of an indirect sort, and if you've read the pandas-related books
# you must have already seen many higher-level examples.
# Given a key or keys (an array of values or multiple arrays of values),
# you wish to obtain an array of integer indices (I refer to them colloquially as indexers)
# that tells you how to reorder the data to be in sorted order.
# Two methods for this are argsort and numpy.lexsort.
# As an example:

values = np.array([5, 0, 1, 3, 2])

values
```

Out[156]:

```
array([5, 0, 1, 3, 2])
```

In [158]:

```
indexer = values.argsort()

indexer
```

Out[158]:

```
array([1, 2, 4, 3, 0], dtype=int64)
```

In [159]:

```
values[indexer]
```

Out[159]:

```
array([0, 1, 2, 3, 5])
```

In [160]:

```
# As a more complicated example, this code reorders a two-dimensional array by its first row:
```

```
arr = np.random.randn(3, 5)
```

```
arr
```

Out[160]:

```
array([[ 0.97544999,  1.63128663, -0.54927156, -0.3058123 ,  0.29766513],
       [ 0.45944142,  1.36198492,  0.49668379, -0.49842716, -0.39958665],
       [-1.2787087 ,  0.81086645, -0.13821823,  0.48103853, -0.44511099]])
```

In [161]:

```
arr[0] = values
```

In [162]:

```
arr
```

Out[162]:

```
array([[ 5.          ,  0.          ,  1.          ,  3.          ,  2.          ],
       [ 0.45944142,  1.36198492,  0.49668379, -0.49842716, -0.39958665],
       [-1.2787087 ,  0.81086645, -0.13821823,  0.48103853, -0.44511099]])
```

In [163]:

```
arr[:, arr[0].argsort()]
```

Out[163]:

```
array([[ 0.          ,  1.          ,  2.          ,  3.          ,  5.          ],
       [ 1.36198492,  0.49668379, -0.39958665, -0.49842716,  0.45944142],
       [ 0.81086645, -0.13821823, -0.44511099,  0.48103853, -1.2787087 ]])
```

In [165]:

```
# Lexsort is similar to argsort, but it performs an indirect lexicographical sort on multiple key arrays.
```

```
# Suppose we wanted to sort some data identified by first and last names:
```

```
first_name = np.array(['Bob', 'Jane', 'Steve', 'Bill', 'Barbara'])
```

```
first_name
```

Out[165]:

```
array(['Bob', 'Jane', 'Steve', 'Bill', 'Barbara'], dtype='<U7')
```

In [166]:

```
last_name = np.array(['Jones', 'Arnold', 'Arnold', 'Jones', 'Walters'])  
last_name
```

Out[166]:

```
array(['Jones', 'Arnold', 'Arnold', 'Jones', 'Walters'], dtype='<U7')
```

In [167]:

```
sorter = np.lexsort((first_name, last_name))  
sorter
```

Out[167]:

```
array([1, 2, 3, 0, 4], dtype=int64)
```

In [168]:

```
zip(last_name[sorter], first_name[sorter])
```

Out[168]:

```
<zip at 0xa01026aac8>
```

In []:

```
# Lexsort can be a bit confusing the first time you use it because the order in which the  
# keys are used to order the data starts with the last array passed.  
# Above, last_name was used before first_name.  
  
# pandas methods like Series's and DataFrame's sort_values method are implemented with variants  
# of these functions (which also must take into account missing values).
```

In [170]:

```
# Alternative Sort Algorithms  
  
# A stable sorting algorithm preserves the relative position of equal elements.  
# This can be especially important in indirect sorts where the relative ordering is meaningful:  
  
values = np.array(['2:first', '2:second', '1:first', '1:second', '1:third'])  
values
```

Out[170]:

```
array(['2:first', '2:second', '1:first', '1:second', '1:third'],  
      dtype='<U8')
```


In [171]:

```
key = np.array([2, 2, 1, 1, 1])  
  
key
```

Out[171]:

```
array([2, 2, 1, 1, 1])
```

In [172]:

```
indexer = key.argsort(kind='mergesort')  
  
indexer
```

Out[172]:

```
array([2, 3, 4, 0, 1], dtype=int64)
```

In [173]:

```
values.take(indexer)
```

Out[173]:

```
array(['1:first', '1:second', '1:third', '2:first', '2:second'],  
      dtype='<U8')
```

In []:

```
# The only stable sort available is mergesort, which has guaranteed  $O(n \log n)$  performance  
# (for complexity buffs), but its performance is on average worse than the default quicksort method.  
# This is not something that most users will ever have to think about,  
# but it's useful to know that it's there.  
  
# A summary of available array sorting methods and their relative performance (and performance guarantees):
```

#	Kind	Speed	Stable	Work space	Worst case
#	'quicksort'	1	No	θ	$O(n^2)$
#	'mergesort'	2	Yes	$n / 2$	$O(n \log n)$
#	'heapsort'	3	No	θ	$O(n \log n)$

< >

In [175]:

```
# Partially Sorting Arrays
```

```
# One of the goals of sorting can be to determine the largest or smallest elements in an array.
```

```
# NumPy has optimized methods, numpy.partition and np.argpartition,  
# for partitioning an array around the k-th smallest element:
```

```
np.random.seed(12345)
```

```
arr = np.random.randn(20)
```

```
arr
```

Out[175]:

```
array([-0.20470766,  0.47894334, -0.51943872, -0.5557303 ,  1.96578057,  
       1.39340583,  0.09290788,  0.28174615,  0.76902257,  1.24643474,  
       1.00718936, -1.29622111,  0.27499163,  0.22891288,  1.35291684,  
       0.88642934, -2.00163731, -0.37184254,  1.66902531, -0.43856974])
```

In [177]:

```
np.partition(arr, 3)
```

Out[177]:

```
array([-2.00163731, -1.29622111, -0.5557303 , -0.51943872, -0.37184254,  
       -0.43856974, -0.20470766,  0.28174615,  0.76902257,  0.47894334,  
       1.00718936,  0.09290788,  0.27499163,  0.22891288,  1.35291684,  
       0.88642934,  1.39340583,  1.96578057,  1.66902531,  1.24643474])
```

In [178]:

```
# After you call partition(arr, 3), the first three elements in the result are the smallest three values
```

```
# in no particular order.
```

```
# numpy.argpartition, similar to numpy.argsort, returns the indices that rearrange the data
```

```
# into the equivalent order:
```

```
indices = np.argpartition(arr, 3)
```

```
indices
```

Out[178]:

```
array([16, 11,  3,  2, 17, 19,  0,  7,  8,  1, 10,  6, 12, 13, 14, 15,  5,  
       4, 18,  9], dtype=int64)
```

In [179]:

```
arr.take(indices)
```

Out[179]:

```
array([-2.00163731, -1.29622111, -0.5557303 , -0.51943872, -0.37184254,
       -0.43856974, -0.20470766,  0.28174615,  0.76902257,  0.47894334,
        1.00718936,  0.09290788,  0.27499163,  0.22891288,  1.35291684,
        0.88642934,  1.39340583,  1.96578057,  1.66902531,  1.24643474])
```

In [181]:

```
# numpy.searchsorted: Finding Elements in a Sorted Array

# searchsorted is an array method that performs a binary search on a sorted array,
# returning the location in the array where the value would need to be inserted to maintain sortedness:

arr = np.array([0, 1, 7, 12, 15])

arr
```

Out[181]:

```
array([ 0,  1,  7, 12, 15])
```

In [182]:

```
arr.searchsorted(9)
```

Out[182]:

```
3
```

In [183]:

```
# You can also pass an array of values to get an array of indices back:

arr.searchsorted([0, 8, 11, 16])
```

Out[183]:

```
array([0, 3, 3, 5], dtype=int64)
```

In [184]:

```
# You might have noticed that searchsorted returned 0 for the 0 element.  
# This is because the default behavior is to return the index at the left side of a group of equal values:
```

```
arr = np.array([0, 0, 0, 1, 1, 1, 1])
```

```
arr
```

Out[184]:

```
array([0, 0, 0, 1, 1, 1, 1])
```

In [185]:

```
arr.searchsorted([0, 1])
```

Out[185]:

```
array([0, 3], dtype=int64)
```

In [186]:

```
arr.searchsorted([0, 1], side='right')
```

Out[186]:

```
array([3, 7], dtype=int64)
```

In [188]:

```
# As another application of searchsorted, suppose we had an array of values between 0 and 10,000,  
# and a separate array of “bucket edges” that we wanted to use to bin the data:
```

```
data = np.floor(np.random.uniform(0, 10000, size=50))
```

```
data
```

Out[188]:

```
array([2449., 7928., 4951., 9150., 9453., 5332., 2524., 7208., 3674.,  
       4986., 2265., 3535., 6508., 3129., 7687., 7818., 8524., 9499.,  
       1073., 9107., 3360., 8263., 8981.,  427., 1957., 2945., 6269.,  
        862., 1429., 5158., 6893., 8566., 6473., 5816., 7111., 2524.,  
       9001., 4422.,  205., 9596., 6522., 5132., 6823., 4895., 9264.,  
       5158.,  721., 5675., 6152., 9415.])
```

In [189]:

```
bins = np.array([0, 100, 1000, 5000, 10000])  
  
bins
```

Out[189]:

```
array([ 0, 100, 1000, 5000, 10000])
```

In [191]:

```
#To then get a labeling of which interval each data point belongs to (where 1 would mean t  
he bucket [0, 100)),  
# we can simply use searchsorted:  
  
labels = bins.searchsorted(data)  
  
labels
```

Out[191]:

```
array([3, 4, 3, 4, 4, 4, 3, 4, 3, 3, 3, 3, 4, 3, 4, 4, 4, 4, 3, 4, 3, 4,  
       4, 2, 3, 3, 4, 2, 3, 4, 4, 4, 4, 4, 4, 3, 4, 3, 2, 4, 4, 4, 4, 3,  
       4, 4, 2, 4, 4, 4], dtype=int64)
```

In [194]:

```
# This, combined with pandas's groupby, can be used to bin data:  
  
pd.Series(data).groupby(labels).mean()
```

Out[194]:

```
2    553.750000  
3   3132.375000  
4   7482.733333  
dtype: float64
```

In [217]:

```
# Writing Fast NumPy Functions with Numba

# Numba is an open source project that creates fast functions for NumPy-like data using CPUs, GPUs,
# or other hardware. It uses the LLVM Project to translate Python code into compiled machine code.

# To introduce Numba, let's consider a pure Python function that computes
# the expression (x - y).mean() using a for loop:

def mean_distance(x, y):
    nx = len(x)
    result = 0.0
    count = 0
    for i in range(nx):
        result += x[i] - y[i]
        count += 1
    return result / count
```

In [199]:

```
#This function is very slow:

x = np.random.randn(10000000)

x
```

Out[199]:

```
array([-1.56565729, -0.56254019, -0.03266414, ..., -2.57718168,
        0.42536042,  1.37299858])
```

In [200]:

```
y = np.random.randn(10000000)

y
```

Out[200]:

```
array([-0.24829772, -0.1873352 , -0.7450831 , ...,  2.10103003,
        0.26415342,  0.73032225])
```

In [203]:

```
%timeit mean_distance(x, y)
```

24 s ± 591 ms per loop (mean ± std. dev. of 7 runs, 1 loop each)

In [204]:

```
%timeit (x - y).mean()
```

231 ms \pm 42.1 ms per loop (mean \pm std. dev. of 7 runs, 1 loop each)

In [218]:

```
# The NumPy version is over 100 times faster.  
# We can turn this function into a compiled Numba function using the numba.jit function:
```

```
import numba as nb
```

```
numba_mean_distance = nb.jit(mean_distance)
```

```
numba_mean_distance
```

Out[218]:

CPUDispatcher(<function mean_distance at 0x000000A01CA935E8>)

In [219]:

```
# We could also have written this as a decorator:
```

```
@nb.jit  
def mean_distance(x, y):  
    nx = len(x)  
    result = 0.0  
    count = 0  
    for i in range(nx):  
        result += x[i] - y[i]  
        count += 1  
    return result / count
```

In [223]:

```
# The resulting function is actually faster than the vectorized NumPy version:
```

```
%timeit numba_mean_distance(x, y)
```

2.04 μ s \pm 168 ns per loop (mean \pm std. dev. of 7 runs, 100000 loops each)

In [224]:

```
# Numba cannot compile arbitrary Python code, but it supports a significant subset of pure Python  
# that is most useful for writing numerical algorithms.  
  
# Numba is a deep library, supporting different kinds of hardware, modes of compilation, and user extensions.  
# It is also able to compile a substantial subset of the NumPy Python API without explicit for loops.  
# Numba is able to recognize constructs that can be compiled to machine code,  
# while substituting calls to the CPython API for functions that it does not know how to compile.  
# Numba's jit function has an option, nopython=True, which restricts allowed code to Python code that can  
# be compiled to LLVM without any Python C API calls. jit(nopython=True) has a shorter alias numba.njit.  
  
# In the previous example, we could have written:
```

```
from numba import float64, njit  
  
@njit(float64(float64[:], float64[:]))  
def mean_distance(x, y):  
    return (x - y).mean()
```

In []:

```
# I encourage you to learn more by reading the online documentation for Numba.
```

In [225]:

```
# Creating Custom numpy.ufunc Objects with Numba  
  
# The numba.vectorize function creates compiled NumPy ufuncs, which behave like built-in ufuncs.  
# Let's consider a Python implementation of numpy.add:  
  
from numba import vectorize  
  
@vectorize  
def nb_add(x, y):  
    return x + y
```


In [226]:

```
# Now we have:
```

```
x = np.arange(10)
```

```
x
```

Out[226]:

```
array([0, 1, 2, 3, 4, 5, 6, 7, 8, 9])
```

In [227]:

```
nb_add(x, x)
```

Out[227]:

```
array([ 0,  2,  4,  6,  8, 10, 12, 14, 16, 18], dtype=int64)
```

In [228]:

```
nb_add.accumulate(x, 0)
```

ValueError Traceback (most recent call last)

<ipython-input-228-e53ce9bd412f> in <module>

----> 1 nb_add.accumulate(x, 0)

ValueError: could not find a matching type for nb_add.accumulate, requested type has type code 'l'

In []:

```
# Advanced Array Input and Output
```

```
# Ordinarily in numpy, np.save and np.load are for storing arrays in binary format on disk.
```

```
# There are a number of additional options to consider for more sophisticated use.
```

```
# In particular, memory maps have the additional benefit of enabling you to work with datasets
```

```
# that do not fit into RAM.
```

In [229]:

```
# Memory-Mapped Files
```

```
# A memory-mapped file is a method for interacting with binary data on disk as though it is stored
```

```
# in an in-memory array.
```

```
# NumPy implements a memmap object that is ndarray-like, enabling small segments of a large file to be read
```

```
# and written without reading the whole array into memory.
```

```
# Additionally, a memmap has the same methods as an in-memory array and thus can be substituted
```

```
# into many algorithms where an ndarray would be expected.
```

```
# To create a new memory map, use the function np.memmap and pass a file path, dtype, shape, and file mode:
```

```
mmap = np.memmap('mymmap', dtype='float64', mode='w+', shape=(10000, 10000))
```

```
mmap
```

Out[229]:

```
memmap([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        ...,
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.]])
```

In [230]:

```
# Slicing a memmap returns views on the data on disk:
```

```
section = mmap[:5]
```

```
section
```

Out[230]:

```
memmap([[0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.],
        [0., 0., 0., ..., 0., 0., 0.]])
```

In [232]:

```
# If you assign data to these, it will be buffered in memory (like a Python file object),  
# but you can write it to disk by calling flush:
```

```
section[:] = np.random.randn(5, 10000)
```

```
section
```

Out[232]:

```
memmap([[ 0.56711132, -0.12717231,  0.52758445, ..., -0.51964959,  
         -0.55576608, -0.62599963],  
        [-0.60118962, -0.2320138 ,  1.78400269, ...,  0.5460359 ,  
         -0.81396878, -0.46026551],  
        [ 0.14608177, -1.1583803 , -1.28189275, ...,  0.53420363,  
         0.09238763, -1.28271782],  
        [-1.5121209 ,  0.9521015 , -0.87179321, ..., -1.28613331,  
         0.01396661, -0.4403933 ],  
        [-0.80654508,  2.11578824, -2.0804114 , ...,  1.12830642,  
         -1.14187654,  0.13829511]])
```

In [233]:

```
mmap.flush()      # writing it to disk
```

In [234]:

```
mmap
```

Out[234]:

```
memmap([[ 0.56711132, -0.12717231,  0.52758445, ..., -0.51964959,  
         -0.55576608, -0.62599963],  
        [-0.60118962, -0.2320138 ,  1.78400269, ...,  0.5460359 ,  
         -0.81396878, -0.46026551],  
        [ 0.14608177, -1.1583803 , -1.28189275, ...,  0.53420363,  
         0.09238763, -1.28271782],  
        ...,  
        [ 0.      ,  0.      ,  0.      , ...,  0.      ,  
         0.      ,  0.      ],  
        [ 0.      ,  0.      ,  0.      , ...,  0.      ,  
         0.      ,  0.      ],  
        [ 0.      ,  0.      ,  0.      , ...,  0.      ,  
         0.      ,  0.      ]])
```

In [235]:

```
del mmap
```

In [236]:

```
# Whenever a memory map falls out of scope and is garbage-collected,  
# any changes will be flushed to disk also.  
# When opening an existing memory map, you still have to specify the dtype and shape,  
# as the file is only a block of binary data with no metadata on disk:
```

```
mmap = np.memmap('mymmap', dtype='float64', shape=(10000, 10000))
```

```
mmap
```

Out[236]:

```
memmap([[ 0.56711132, -0.12717231,  0.52758445, ..., -0.51964959,  
         -0.55576608, -0.62599963],  
        [-0.60118962, -0.2320138 ,  1.78400269, ...,  0.5460359 ,  
         -0.81396878, -0.46026551],  
        [ 0.14608177, -1.1583803 , -1.28189275, ...,  0.53420363,  
         0.09238763, -1.28271782],  
        ...,  
        [ 0.          ,  0.          ,  0.          , ...,  0.          ,  
         0.          ,  0.          ],  
        [ 0.          ,  0.          ,  0.          , ...,  0.          ,  
         0.          ,  0.          ],  
        [ 0.          ,  0.          ,  0.          , ...,  0.          ,  
         0.          ,  0.          ]])
```

In []:

```
# Memory maps also work with structured or nested dtypes as described in a previously
```

In []:

```
# Performance Tips
```

```
# Getting good performance out of code utilizing NumPy is often straightforward,  
# as array operations typically replace otherwise comparatively extremely slow pure Python loops.
```

```
# The following list briefly summarizes some things to keep in mind:
```

```
# • Convert Python loops and conditional logic to array operations and boolean array operations  
# • Use broadcasting whenever possible  
# • Use arrays views (slicing) to avoid copying data  
# • Utilize ufuncs and ufunc methods
```

```
# If you can't get the performance you require after exhausting the capabilities provided by NumPy alone,
```

```
# consider writing code in C, Fortran, or Cython.
```

In [237]:

```
# The Importance of Contiguous Memory

# In some applications the memory layout of an array can significantly affect the speed of
computations.
    # This is based partly on performance differences having to do with the cache hierarch
y of the CPU;
    # operations accessing contiguous blocks of memory (e.g., summing the rows of a C orde
r array)
    # will generally be the fastest because the memory subsystem will buffer the appropria
te blocks of memory
    # into the ultrafast L1 or L2 CPU cache.
    # Also, certain code paths inside NumPy's C codebase have been optimized for the conti
guous case
    # in which generic strided memory access can be avoided.

# To say that an array's memory layout is contiguous means that the elements are stored in
memory in the order
    # that they appear in the array with respect to Fortran (columnmajor) or C (row major)
ordering.
    # By default, NumPy arrays are created as Ccontiguous or just simply contiguous.
    # A column major array, such as the transpose of a C-contiguous array,
    # is thus said to be Fortran-contiguous.
    # These properties can be explicitly checked via the flags attribute on the ndarray:

arr_c = np.ones((1000, 1000), order='C')

arr_c
```

Out[237]:

```
array([[1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       ...,
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.]])
```

In [238]:

```
arr_f = np.ones((1000, 1000), order='F')

arr_f
```

Out[238]:

```
array([[1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       ...,
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.],
       [1., 1., 1., ..., 1., 1., 1.]])
```

In [239]:

```
arr_c.flags
```

Out[239]:

```
C_CONTIGUOUS : True
F_CONTIGUOUS : False
OWNDATA : True
WRITEABLE : True
ALIGNED : True
WRITEBACKIFCOPY : False
UPDATEIFCOPY : False
```

In [240]:

```
arr_f.flags
```

Out[240]:

```
C_CONTIGUOUS : False
F_CONTIGUOUS : True
OWNDATA : True
WRITEABLE : True
ALIGNED : True
WRITEBACKIFCOPY : False
UPDATEIFCOPY : False
```

In [242]:

```
arr_c.flags.c_contiguous
```

Out[242]:

```
True
```

In [243]:

```
arr_c.flags.f_contiguous
```

Out[243]:

```
False
```

In [244]:

```
arr_f.flags.c_contiguous
```

Out[244]:

```
False
```

In [245]:

```
arr_f.flags.f_contiguous
```

Out[245]:

```
True
```

In [250]:

```
# In this example, summing the rows of these arrays should, in theory, be faster for arr_c  
than arr_f  
# since the rows are contiguous in memory.  
  
# Here I check for sure using %timeit:  
  
%timeit arr_c.sum(1)
```

5.77 ms \pm 782 μ s per loop (mean \pm std. dev. of 7 runs, 100 loops each)

In [251]:

```
%timeit arr_f.sum(1)
```

2.54 ms \pm 944 μ s per loop (mean \pm std. dev. of 7 runs, 100 loops each)

In [252]:

```
# When you're looking to squeeze more performance out of NumPy, this is often a place to i  
nvest some effort.  
# If you have an array that does not have the desired memory order,  
# you can use copy and pass either 'C' or 'F':  
  
arr_f.copy('C').flags
```

Out[252]:

```
C_CONTIGUOUS : True  
F_CONTIGUOUS : False  
OWNDATA : True  
WRITEABLE : True  
ALIGNED : True  
WRITEBACKIFCOPY : False  
UPDATEIFCOPY : False
```

In [253]:

```
arr_f.copy('F').flags
```

Out[253]:

```
C_CONTIGUOUS : False  
F_CONTIGUOUS : True  
OWNDATA : True  
WRITEABLE : True  
ALIGNED : True  
WRITEBACKIFCOPY : False  
UPDATEIFCOPY : False
```

In [254]:

```
arr_c.copy('F').flags
```

Out[254]:

```
C_CONTIGUOUS : False
F_CONTIGUOUS : True
OWNDATA : True
WRITEABLE : True
ALIGNED : True
WRITEBACKIFCOPY : False
UPDATEIFCOPY : False
```

In [255]:

```
arr_c.copy('C').flags
```

Out[255]:

```
C_CONTIGUOUS : True
F_CONTIGUOUS : False
OWNDATA : True
WRITEABLE : True
ALIGNED : True
WRITEBACKIFCOPY : False
UPDATEIFCOPY : False
```

In [256]:

```
# When constructing a view on an array, keep in mind that the result is not guaranteed to be contiguous:
```

```
arr_c[:50].flags.contiguous
```

Out[256]:

```
True
```

In [257]:

```
arr_c[:, :50].flags
```

Out[257]:

```
C_CONTIGUOUS : False
F_CONTIGUOUS : False
OWNDATA : False
WRITEABLE : True
ALIGNED : True
WRITEBACKIFCOPY : False
UPDATEIFCOPY : False
```

The End