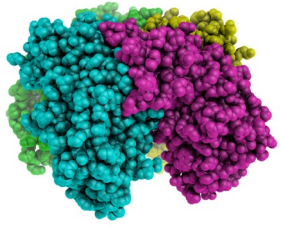


Eines Informatiques: Python

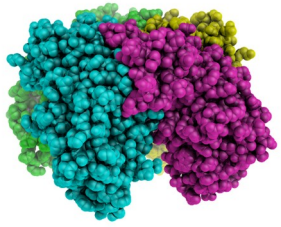
Ramon Crehuet

Curs 2020-2021

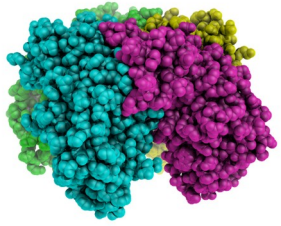


Overview

- Why Python
- Language basics
- Functions and modules
- Working with files
- Classes and objects (bare minimum!)
- Working with arrays: Numpy
- Data visualization
- Scientific modules. Scipy
- Other scientific modules: Pandas, scikit-learn, biopython
- Profiling and optimization
- Beyond Python

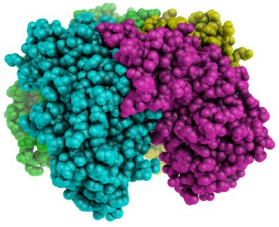


Introduction



Overview

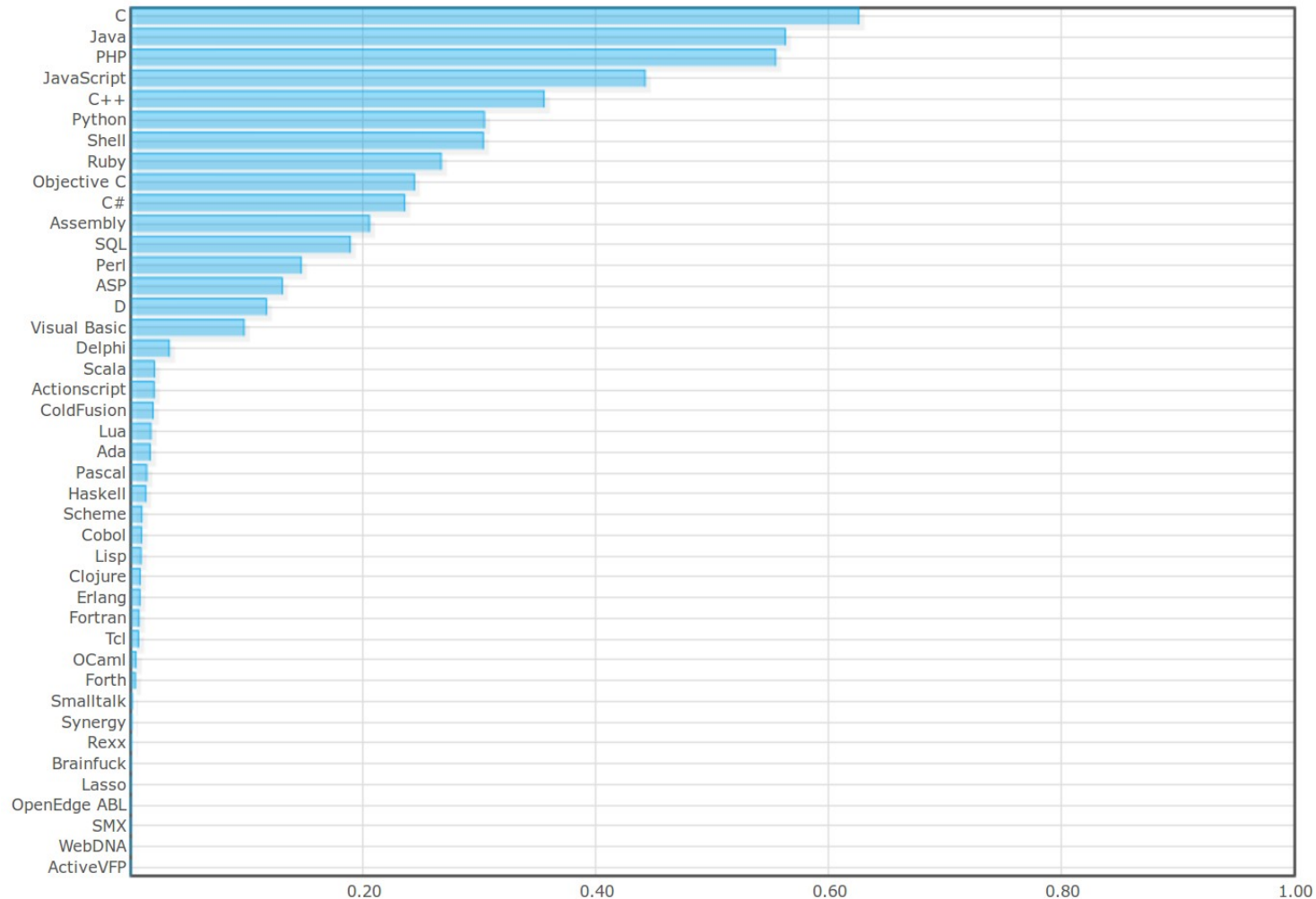
- Why Python
- Language basics
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- Data visualization
- Scientific modules. Scipy
- Organizing data: pandas and xarray
- Other scientific modules: scikit-learn, biopython
- Profiling and optimization
- Beyond Python



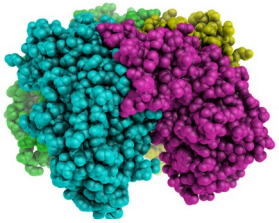
Language popularity

Normalized Comparison

This is a chart showing combined results from all data sets, listed individually below.



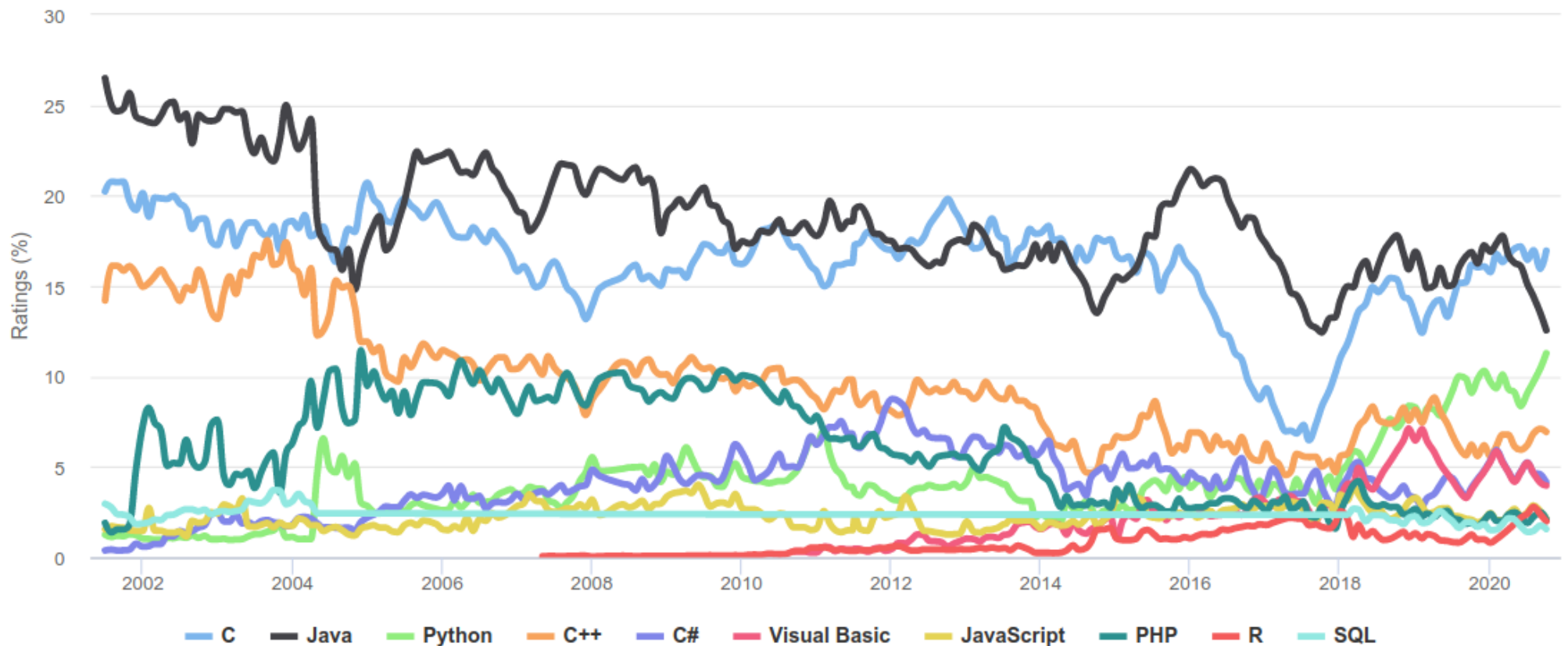
<http://langpop.com/>



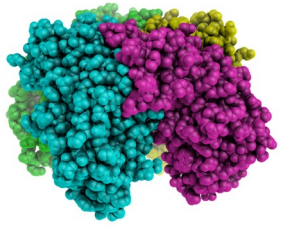
Language popularity

TIOBE Programming Community Index

Source: www.tiobe.com

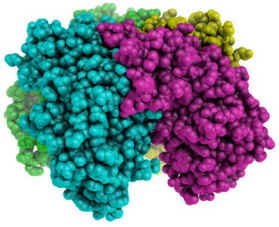


<http://www.tiobe.com/index.php/content/paperinfo/tpci/index.html>



Python for science

- A high level language gives more time to more complex problems
 - At the expense of hiding important details
- Example:
 - A reaction mechanism
 - Optimisation of an energy function
 - Steepest descent, conjugate gradients, quasi-Newton
 - Implementation of BFGS quasi-Newton
 - Memory issues, diagonalization, matrix inversion...
 - Calculation of numerical gradients or hessians:
 - machine precision, central differences, etc.



Python for science

Compiled languages

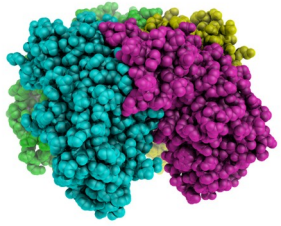
Fast
Difficult
non-interactive

Matlab, Mathematica, Octave

Slow
Rich libraries
Nice development environment
Restricted base language
Expensive (some)

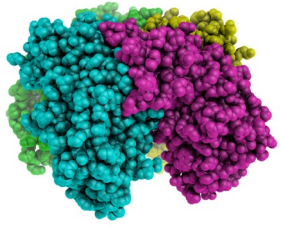
Python

Rich libraries (less than matlab)
Other libraries
Free
Active community
Harder than Matlab



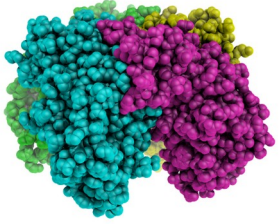
Matlab, Mathematica?

- Scientific computing:
 - ipython + scipy + matplotlib
- Free
- Open source
- Extensible
- Bioinformatics
 - Biopython
- Molecular Dynamics
 - MMTK
- Efficiency
 - Numba, Cython, Cupy, Fortran, C
- Server control
- XML parser



Low level vs. high level

- Python is a high level language
- You can focus on:
 - Low level issues
 - Higher complexity of problems
- Low level issues
 - Variable types
 - Machine precision
- But also
 - Extend
 - Mantain
 - Document code



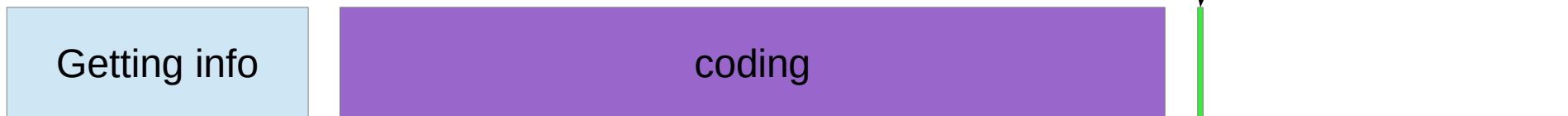
Python vs. Fortran

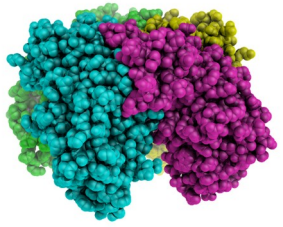
Different time distribution to get a task done

Python

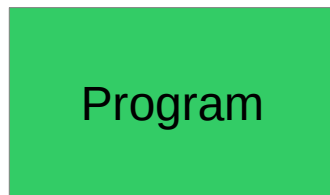
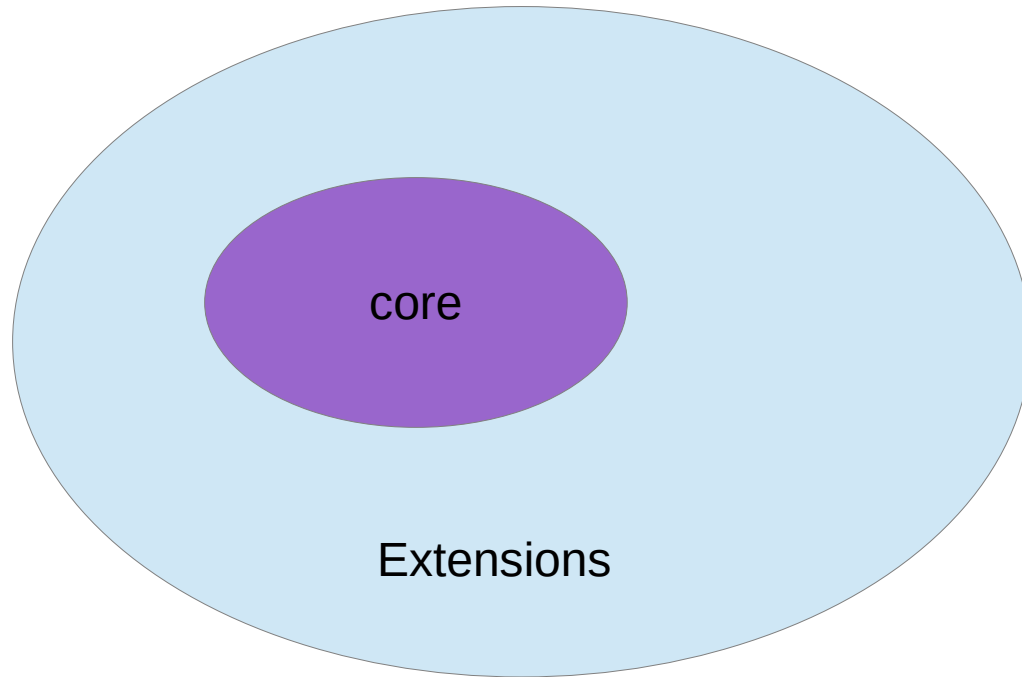


Fortran

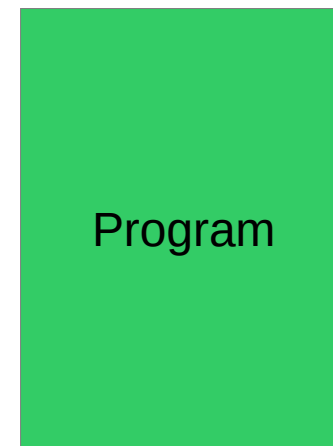
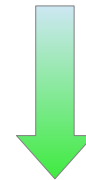
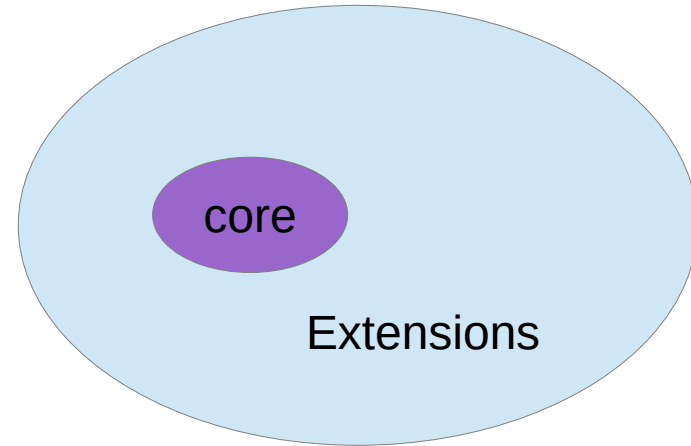


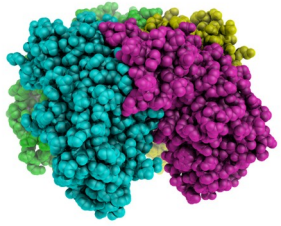


Python



Fortran





Python for science

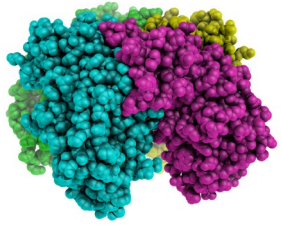
- The homogenization of scientific computing, or why Python is steadily eating other languages' lunch

<http://www.talyarkoni.org/blog/2013/11/18/the-homogenization-of-scientific-computing-or-why-python-is-steadily-eating-other-languages-lunch/>

- 10 Reasons Python Rocks for Research (And a Few Reasons it Doesn't)

<http://www.stat.washington.edu/~hoytak/blog/whypython.html>

-

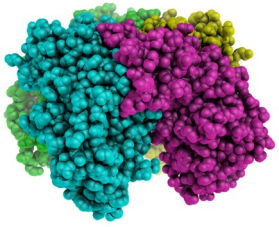


Hello World program

```
print("Hello World!")
```

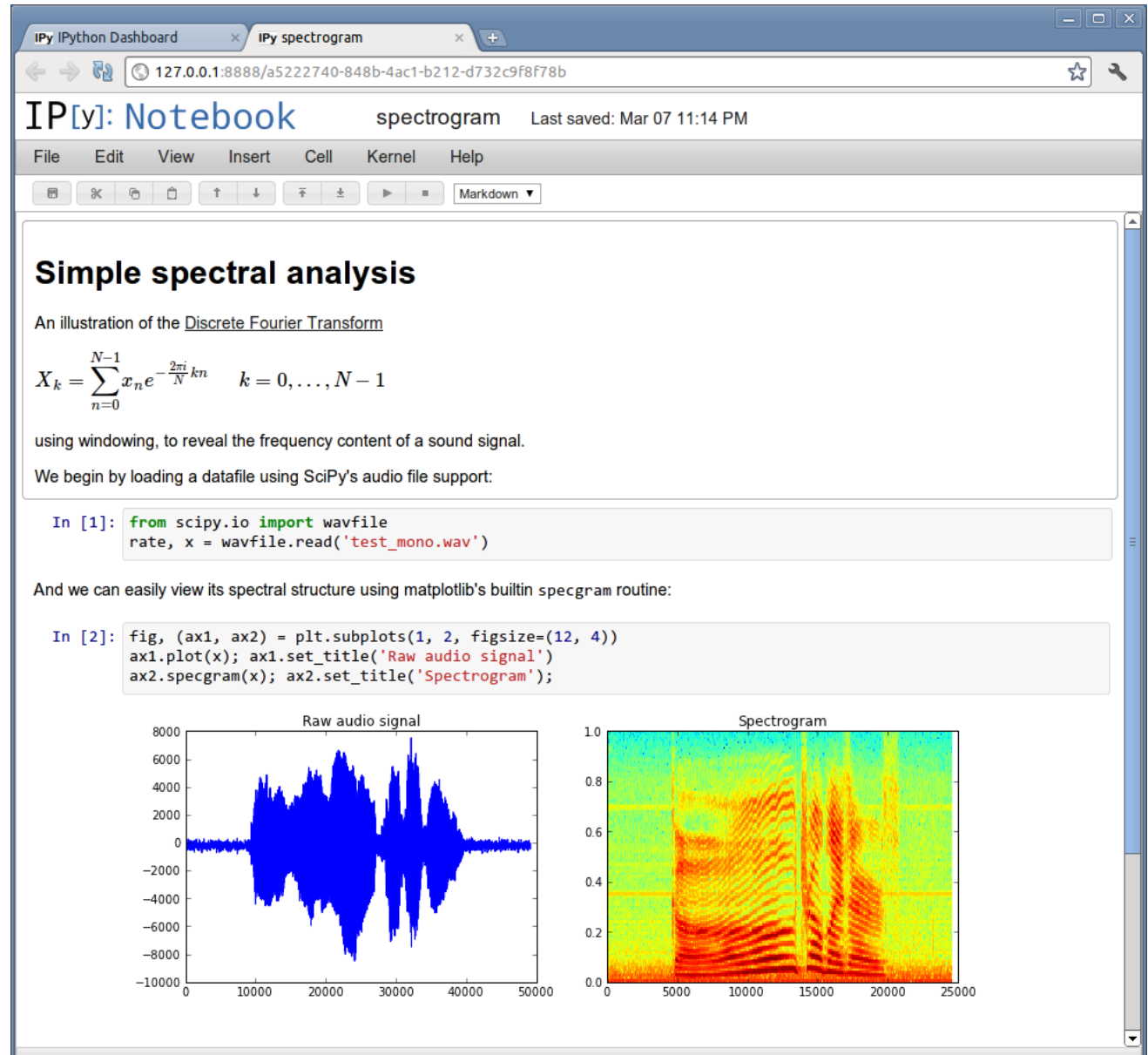
```
print("Hello World!")
```

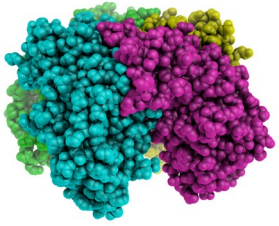
```
$ python3 hello.py
```



Interactive shells

- python
- IDLE
- **ipython**
 - shell
 - notebook
- spyder
- visual studio code
- PIDA
- Sage





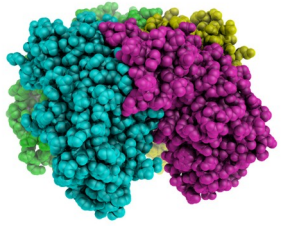
Dynamically typed

```
>>> a = 4
>>> type(a)
<class 'int'>
>>> b = 7.6
>>> type(b)
<class 'float'>
>>> type(a+b)
<class 'float'>
>>> c = 'Hola'
>>> c + ' Que tal?'
'Hola Que tal?'
>>> c + a
```

Traceback (most recent call last):

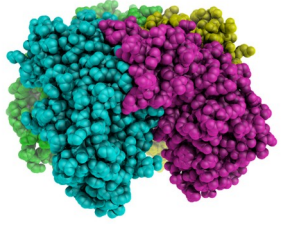
File "<stdin>", line 1, in <module>

TypeError: Can't convert 'int' object to str implicitly

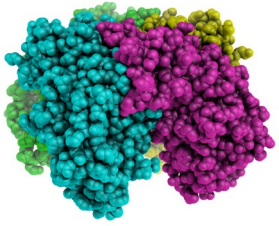


Which python version?

- Language is fast evolving
- 2 versions now coexist: 3.x and 2.x
- These versions are not completely compatible
- 3.x is better and continued
- 2.x has some software still not ported
- Both can safely coexist
 - Packages and shells are for a specific version
- `2to3 -w hello.py`



Language elements



Numbers

Integers:

```
> i = 5  
> j = i**i**i
```

Limited by amount of memory:

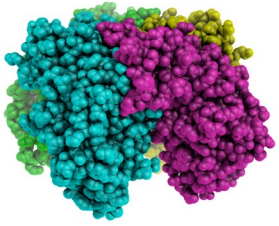
```
>>> i.bit_length()  
3  
  
>>> j = i**i**i  
>>> j.bit_length()  
7257  
  
>>> 9 % 5 #modulo  
4
```

Floating point:

```
>>> x = 5.  
>>> y = x**x**x  
Traceback (most recent call last):  
  File "<stdin>", line 1, in <module>  
OverflowError: (34, 'Numerical result  
out of range')
```

Division vs integer division (Python 3):

```
>>> 3/2  
1.5  
  
>>> 3//2  
1  
  
>>> j/i #Returns a Float
```



Assignments

Explicit notation:

```
> i = i+1  
> j = j / 10.
```

Short notation:

```
> i += 1  
> j /= 10.
```

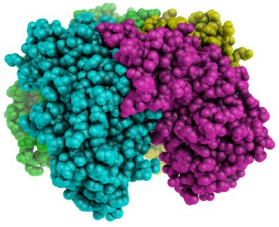
Floating point:

```
>>> x = 5.  
>>> y = x**x**x
```

Traceback (most recent call
last):

File "<stdin>", line 1, in
<module>

OverflowError: (34, 'Numerical
result out of range')



strings

Strings:

```
> str(6.7)
> c = 'Hola'
```

Operations:

```
> s = 'numeric ' + 'python'
> len(s)
> s[5]
'i'
> s.split()
['numeric', 'python']
```

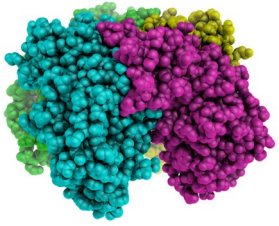
```
> print('Result: %5.3f' % (11./3.))
3.667
```

Non mutable:

```
> s[6]
> s[6]='7'
```

Regular expressions

```
import re
```



Lists, sets and tuples

- Fortran **dimension**:

much more flexible

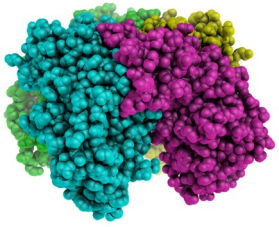
```
> l=[6, 'a', [5,[9,8,7,6]], -  
    6.5, (True, True)]  
> [1,2]+[3,4]  
> l.append(6)
```

- sets:

```
s=set([4,3,2,3])  
> 4 in s  
True  
> s  
set([2, 3, 4])
```

- Tuples are unmutable lists

```
t=(1,2,3)
```



Lists, sets and tuples

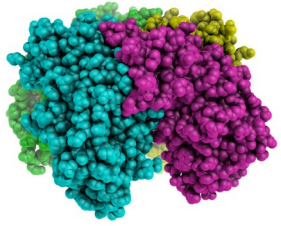
List indexing and methods:

```
> l = list(range(10))  
> l[4] = 20  
> l[4:]  
> l[-4]  
> l[:]  
> l[::-1] #reverse  
> l.reverse()  
> l.pop()  
> l.extend([3,4,5])  
> l.sort()
```

First
index is 0

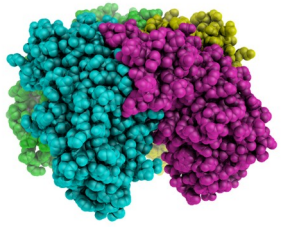
Set methods:

```
> s1=set([1,2,3,4])  
> s2=set([3,4,5,6,7])  
> s1.union(s2)  
> s1.intersection(s2)  
> s1.difference(s2)  
> s2.difference(s1)  
> s1.intersection(s2) == s2 & s1  
True  
> s1 - s2 == s2 - s1  
False
```



Uses of lists, sets and tuples

- Calculate and keep all the primes < 1000
- Given a coordinate file, calculate for each atom a list of all the atoms that are at less than 0.2nm.
- Get the solutions of a quadratic equation (0,1,2) or (real vs. complex).
- <http://docs.python.org/3/tutorial/datastructures.html>



Copying and looping over lists

lists are treated as pointers:

copying lists, makes a copy of the pointer.

```
> l=[1,2,3,4]
```

```
> l2=l
```

```
> l[2]=1000
```

```
l1
```

```
[1, 2 , 1000, 4]
```

Looping over lists:

Fortran/C style:

```
num=[2,3,2,3,4,5,5]
```

```
for i in range(len(num)):
```

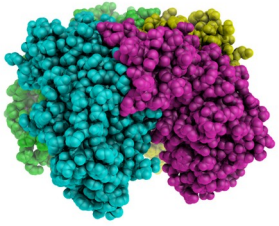
```
    print(num[i])
```

Pythonic style:

```
for item in num:
```

```
    print(item)
```

This can be used for sets,
dictionaries, and tuples.



Dictionaries

Setting elements:

```
> phone={}
> phone['Ramon']='1242'
> phone['Joan']='1323'
> phone['Quique']='1242'
> phone.keys()
['Quique', 'Joan', 'Ramon']
> d2 = dict(Ramon=1242, Joan=1323,
            Quique=1242)
```

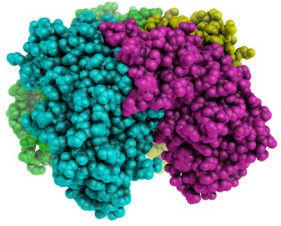
Dictionaries are not ordered

Getting elements:

```
> for key in phone:
...     print(key, phone[key])
Quique 1242
Joan 1323
Ramon 1242
```

Removing elements:

```
> del(phone['Ramon'])
```



The beauty of Python blocks

We are usually told to indent blocks for clarity.

Python makes this the syntax rule to identify blocks.

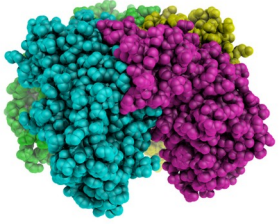
The code has to be nice!

Convention:

- Use 4 spaces
- Use spaces, not tabs.

```
while iter < maxIter:  
    x = f(x)  
    iter = iter + 1
```

```
if i>0:  
    print("i is positive")  
elif i==0:  
    print("i is zero")  
else:  
    print("i is negative")
```



Execution control: if

if... elif... else

```
if <condition>:
```

```
    <block>
```

```
elif <condition>:
```

```
    <block>
```

```
else:
```

```
    <block>
```

```
4==4 #True
```

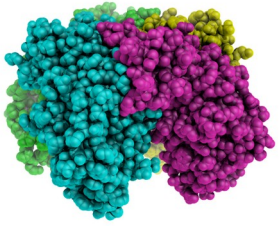
```
5!=4 #True
```

```
4>=5 #False
```

```
4 in [4,5] #True
```

```
result=True
```

```
if result: print('yes')
```



Execution control

Conditions can be combined with:

and or not ()

Object identity:

```
> a=[1,2,3]
```

```
> b=a
```

```
> b is a
```

```
True
```

Any non-zero number or non-empty string is True:

```
> if []: print ('yes')  
      else: print('no')
```

```
no
```

```
> if 5 and 'result':  
      print('yes')
```

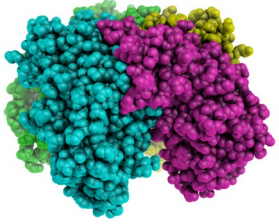
```
else:
```

```
      print('no')
```

```
yes
```

```
> if 5 or 1/0: print('yes')
```

```
yes
```



for and while loops

For loops

```
> dict={4:'a',3:'b', 2:'c',  
        1:'d'}  
> for i in dict:  
    print(i, dict[i])
```

While

```
while <condition>:  
    <block>
```

Break continue pass

```
> pass # does nothing
```

break: Fortran EXIT

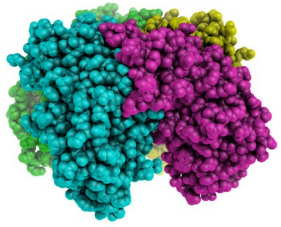
```
if x>0:
```

```
    pass
```

```
else:
```

```
    break
```

cycle: Fortran CONTINUE



list comprehension and enumerate

simple way to create lists:

```
> l=[x**2 for x in range(8)]
```

```
[0, 1, 4, 9, 16, 25, 36, 49]
```

with conditionals:

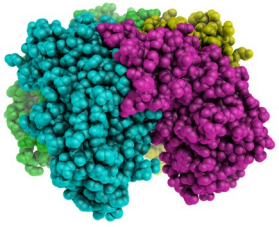
```
l2= [(i, -2*i+3) for i in l if i % 3 == 0]
```

```
[(0, 3), (9, -15), (36, -69)]
```

Nested lists:

```
> [(x, y) for x in [1,2,3] for y in [3,1,4] if x != y]
```

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



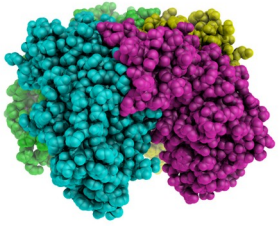
list comprehension and enumerate

Enumerate indexes lists:

```
line='how do you do?'  
line=line.split()  
for i, word in enumerate(line):  
    print(i, word.upper())  
0 HOW  
1 DO  
2 YOU  
3 DO?
```

Enumerate returns an iterator

```
> enumerate(['a', 'b', 'c'])  
<enumerate object at 0x1ebeaa50>
```

Be pythonic

Convert the negative elements of a list to positive

```
>>> x = [1, 2, -4, -5, 3, -5]
```

```
j = 0
```

```
while j < len(x):
```

```
    x[j] = abs(x[j])
```

```
    j += 1
```

```
for j in range(len(x)):
```

```
    x[j] = abs(x[j])
```

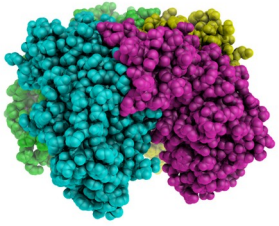
Or with list comprehensions

```
x = [abs(j) for j in x]
```

Or with functional programming

```
x= map(abs, x) #returns an iterator
```

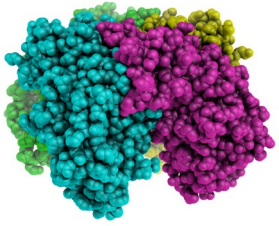
<http://docs.python-guide.org/en/latest/writing/style/>



More python functions

```
print(3,4,5, sep='o', end='<<<\n')  
zip([1,2,3], ['a', 'b', 'c', 'd'])  
a = input('Write a number: ')  
len([1,2,3])  
list(range(5))  
range(20,10,-1)  
sorted([5,4,3,5])  
sum([5,4,3,5])
```

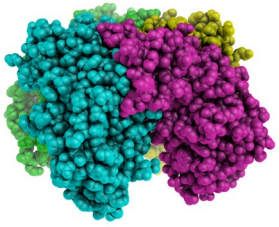
<http://docs.python.org/3.3/library/functions.html>



Mutable and immutable

- Mutable objects can be mutated.
 - Their identity remains the same
- Immutable objects are “mutated” by creating a new object

```
>>> a = 4
>>> id(a)
9157088
>>> a += 2
>>> id(a)
9157152
>>> s = 'Hola'
>>> id(s)
140165884365656
>>> s = s + ' que tal?'
>>> id(s)
140165884365712
>>> ll = [3,4,5]
>>> id(ll)
140165884674416
>>> ll.append(6)
>>> id(ll)
140165884674416
```



Identity and equality

```
>>> 1.0 is 1.0
```

```
True
```

```
>>> 1.0 == 1.0
```

```
True
```

```
>>> 1 == 1.0
```

```
True
```

```
>>> 1 is 1.0
```

```
False
```

```
>>> a = 4
```

```
>>> b = a
```

```
>>> a is b
```

```
True
```

```
>>> id(a)
```

```
9157088
```

```
>>> id(b)
```

```
9157088
```

```
>>> l1 = [1,2,3,]
```

```
>>> l2 = l1
```

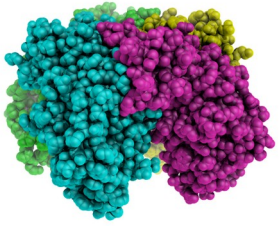
```
>>> l2 = l1[:]
```

```
>>> l2 is l1
```

```
False
```

```
>>> l2 == l1
```

```
True
```



Objects: everything

```
>>> a = 5
>>> isinstance(a, int)
True
>>> object
<class 'object'>
>>> int
<class 'int'>
>>> isinstance(a, object)
True
>>> issubclass(int, object)
True
```

Objects have variables:

```
> c = 4+5j
```

```
> c.real
```

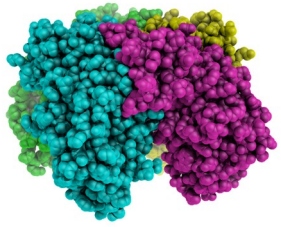
Objects have methods:

```
> c.conjugate #the method
```

```
> c.conjugate() #its call
```

And we can apply functions to objects:

```
> abs(c)
```



Python flow with pythontutor

www.pythontutor.com/visualize.html#mode=display

Most Visited ▾ Macromolecular M... ▾ Portal de Acceso a la ... ▾ Computational Che... ▾ Save to Mendeley

```

1 x = [1, 2, 3]
2 y = [4, 5, 6]
3 z = y
4 y = x
5 x = z
6
7 x = [1, 2, 3] # a different [1, 2, 3] list!
8 y = x
9 x.append(4)
10 y.append(5)
11 z = [1, 2, 3, 4, 5] # a different list!
12 x.append(6)
13 y.append(7)
14 y = "hello"
15
16
17 def foo(lst):
18     lst.append("hello")

```

[Edit code](#)

Frames

Global frame	
x	[1, 2, 3]
y	"hello"
z	[1, 2, 3, 4, 5]
foo	
bar	

Objects

list	
0	1
1	2
2	3
3	4
4	5
5	6
6	7
7	"hello"

list	
0	1
1	2
2	3
3	4
4	5
5	"hello"

function
foo(lst)

function
bar(myLst)

Program terminated

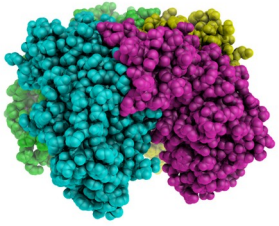
→ line that has just executed
→ next line to execute

Program output:

```

[1, 2, 3, 4, 5, 6, 7, 'hello']
[1, 2, 3, 4, 5, 'hello']

```



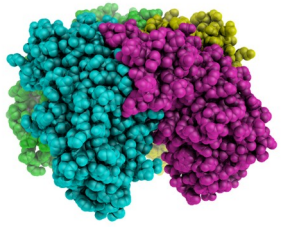
try... except

“Look before you leap”:

```
def safe_divide_1(x, y):  
    if y==0:  
        print("Divide-by-0 attempt  
detected")  
        return None  
    else:  
        return x/y
```

“It's easier to ask forgiveness
than permission”:

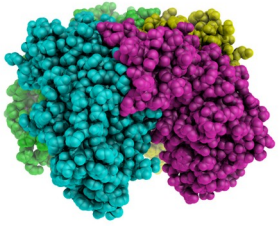
```
def safe_divide_2(x, y):  
    try:  
        return x/y  
    except ZeroDivisionError:  
        print("Divide-by-0 attempt  
detected")  
        return None
```



Short notebook tutorial

But watch “I don’t like notebooks”:

<http://ipython.org/ipython-doc/dev/interactive/tutorial.html>



beyond python

TAB autocomplete:

- functions
- methods
- files
- ...

reload command

cursor keys get history:

- even previous sessions!
- text + keys: previous match

?: intro to ipython

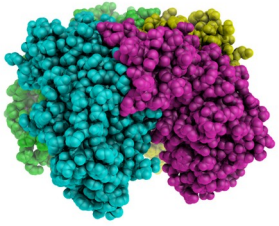
%quickref

Ctrl-r : previous commands

Without ipython:

`python3 -u script.py` enters interactive mode

```
>>> import rlcompleter, readline
>>> readline.parse_and_bind('tab:complete')
```



Magic functions

`%timeit x=10` : time the 'x=10' statement with high precision.

`%%timeit x=2**100`

`x*100` : time 'x*100' with a setup of 'x=2**100'; setup code is not counted. This is an example of a cell magic.

`%cpaste, %paste`: Paste & execute a pre-formatted code block from clipboard.

`%history`

`%load_ext`

`%run`

`%pdb`: Control the automatic calling of the pdb interactive debugger.

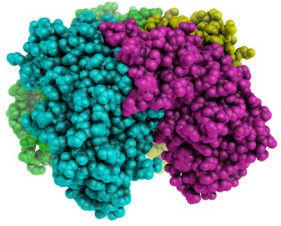
`%pylab`

`%timeit`

`%pwd`

`%cd`

`%%bash` <http://ipython.org/ipython-doc/dev/interactive/tutorial.html>



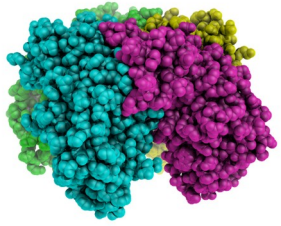
running scripts

`%run script.py`

`import script.py`

are not the same!

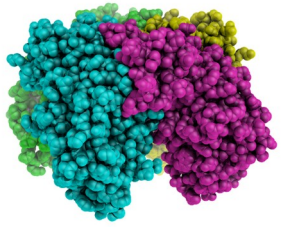
`%run script.py` is like `python3 script.py`



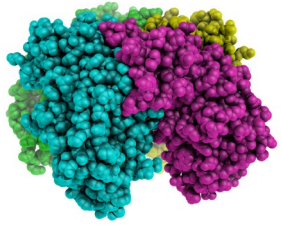
ipython notebook

- Nice presentation
- Allows parallel execution
- Combines text and code
- Executable or exportable to:
 - html
 - LaTeX
 - python
- Start with: `ipython3 notebook`
- Examples:

<https://github.com/jrjohansson/scientific-python-lectures>

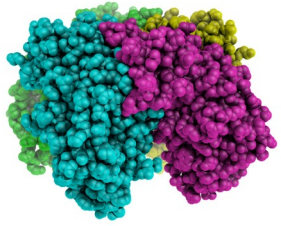


Files



Files

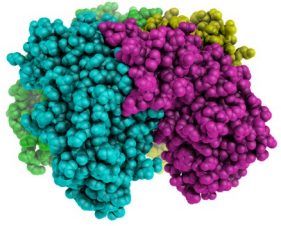
- Files can be text or binary files
- Files can be opened for read, write or append
 - 'r', 'w', 'a+'
- with `open('name') as filein:`
 - Allows automatic file closure



Reading / Writing Files

```
file_in=open('indata.txt','r')  
file_out=open('outdata.txt','w')  
for line in file_in:  
    # Take some information (split() method is very useful!)  
    x = float(line.split()[0])  
    # Apply a given function (fact)  
    fx = fact(x)  
    # Write the result in an output file with a defined format  
    file_out.write('{%:010.3f}\n'.format(fx))
```

But for loading numerical data Numpy is more efficient...



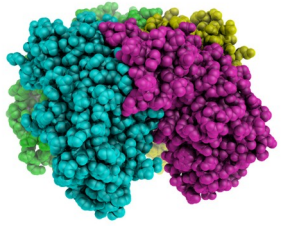
File parsing

- The basic:

```
for line in filein:  
    do something
```

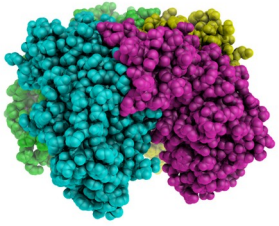
- Common things:

```
if 'optimized' in line: do something  
line = line.split()  
if line.upper().startswith('GEOM'): ...  
energy = float(line[2])
```

skipping lines

- Lines can be skipped by calling `next()` to a file:
for line in filein:
 if 'Optimized' in line:
 next(filein); next(filein) #skip two lines
 do something...



Formatting

- There are several function:

```
'12'.rjust(5), '12'.zfill(5)
```

- But format is more general:

```
print('{0:2d} {1:3d}'.format(x, x*x))
```

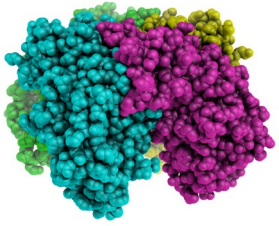
```
print("{:10.3f} {:10.3f} {:10.3f}".format(x,y,z))
```

- List of unknown length:

```
vals = np.linspace(0,1,11)
```

```
print((len(vals)*"{:10.2e} ").format(*vals))
```

<http://docs.python.org/3/library/string.html#formatspec>



Useful modules

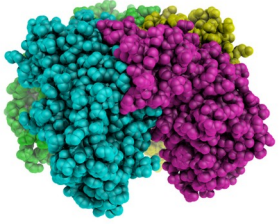
- Similar to `ls`:

```
import glob
files = glob.glob(pattern)
```

- Working with shell-like commands:

```
import os
os.rename(src, dst)
os.mkdir(path)
os.chown(path, uid, gid)
os.getenv(key)
os.walk(directory)
```

<http://docs.python.org/3/library/os.html>



Useful modules

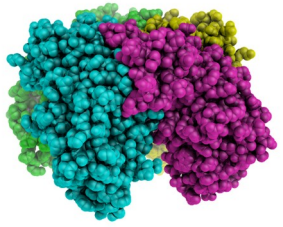
- Reading Excel files:

```
import xlrd
```

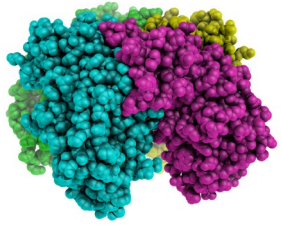
- Working with image files:

```
from PIL import Image
```

<http://www.python-excel.org/>
<http://pillow.readthedocs.org/en/latest/>



Numpy



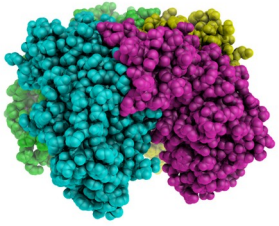
Why Numpy / Scipy?

- Python (alone) is not efficient for numerical calculations
- Python (alone) is not practical for array manipulation
- Numpy provides the data types and methods for arrays
- Scipy provides more elaborate numerical methods
 - Optimization
 - Fast Fourier Transform
 - Linear algebra, etc

```
import numpy as np
```

```
import scipy.optimization
```

```
import scipy.stats as stats
```



numpy arrays

- without numpy:

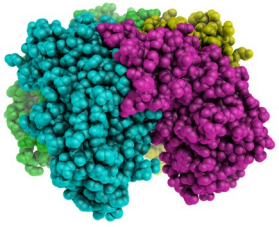
```
> a=[[1,2],[3,4]]  
> b=[[10,20], [30,40]]  
> a+b  
[[1, 2], [3, 4], [10, 20], [30, 40]]
```

- with numpy:

```
> a=np.array(a)  
> b=np.array(b)  
> a+b  
array([[11, 22],[33, 44]])
```

- Array creation

```
a=np.array([1,2,3,4]).reshape([2,2])  
a=np.array([[1,2], [3,4]])  
a=np.zeros([2,2], dtype=int)  
a[0,0]=1.  
a=np.ones((4,4))  
a=np.arange(10)  
a=np.diag([1,2,3,4])  
a=np.tile(a, (10,2))  
a=np.identity(3)  
a=np.linspace(-5,5, 20)
```



Ufuncs

- Unary:

`a.min()`

`a.sum()`

`a.cumsum()`

`a.mean()`

`np.argmin(a)`

`np.exp(-a)`

`np.cov(a)`

`a.tolist()`

- Binary:

`a + b`

`np.dot(a, b)`

- Applying to parts of an array:

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

```
> a.min(axis=0)
```

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

```
a.sum(axis=1)
```

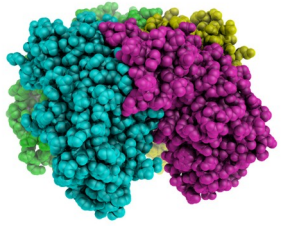
```
array([3, 7])
```

- Python functions are less efficient than numpy functions:

`a.sum()` better than `sum(a)`

`np.min(a)` better than `min(a)`

many implemented as methods and functions



Accessing array elements

- Slicing:

```
> a[2:5]
```

```
> b[:, ::5]
```

```
> a[1:4, ...]
```

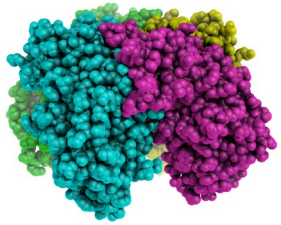
- Fancy indexing:

- Boolean arrays (masks):

```
> a = np.arange(10,15)
> indices = (a**2 > 115) & (a < 14)
> a[indices]
array([11, 12, 13])
```

- With lists:

```
> a = np.arange(10,15)
> y=a[[4,4,1]]
> y
array([14, 14, 11])
> a[[4,4,1]] = [-2, -4, 5]
> a
array([10,  5, 12, 13, -4])
```



Accessing array elements

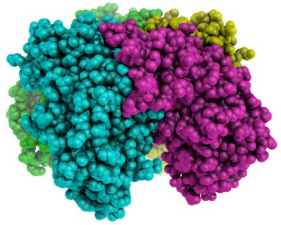
```
>>> a[0,3:5]  
array([3,4])
```

```
>>> a[4:,4:]  
array([[44, 45],  
       [54, 55]])
```

```
>>> a[:,2]  
array([2,12,22,32,42,52])
```

```
>>> a[2::2,::2]  
array([[20,22,24]  
       [40,42,44]])
```

0	1	2	3	4	5
10	11	12	13	14	15
20	21	22	23	24	25
30	31	32	33	34	35
40	41	42	43	44	45
50	51	52	53	54	55



Accessing array elements

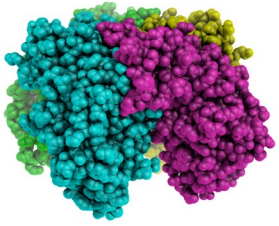
```
>>> a[(0,1,2,3,4),(1,2,3,4,5)]  
array([ 1, 12, 23, 34, 45])
```

```
>>> a[3:,[0, 2, 5]]  
array([[30, 32, 35],  
       [40, 42, 45]],  
      [50, 52, 55])
```

```
>>> mask = array([1,0,1,0,0,1],  
                 dtype=bool)
```

```
>>> a[mask,2]  
array([2,22,52])
```

0	1	2	3	4	5
10	11	12	13	14	15
20	21	22	23	24	25
30	31	32	33	34	35
40	41	42	43	44	45
50	51	52	53	54	55



Accessing array elements

- Slices return views

```
> a = np.arange(5)
> y=a[2:5]
> y *= -1
> a
array([ 0,  1, -2, -3, -4])
> y.flags.owndata
False
```

- np.where

```
> np.where((a>=2)&(a<4), a**2, -1)
```

```
Array([-1, -1,  4,  9, -1])
```

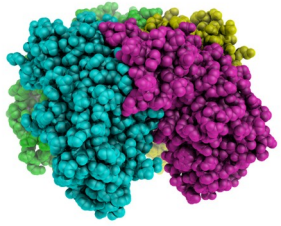
- np.choose
 - Powerful, but complex!
- np.nonzero

- Boolean arrays return copies

```
> a = np.arange(5)
> y = a[a>1-5]
> y *= -1
> a
array([0, 1, 2, 3, 4])
> y.flags.owndata
True
```

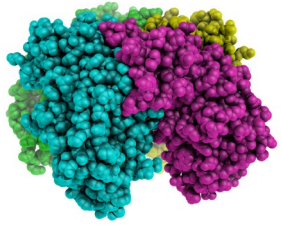
- Fancy indexing returns copies:

```
> a = np.arange(5)
> y=a[[2,3,4]]
> y *= -1
> a
array([0, 1, 2, 3, 4])
> y.flags.owndata
True
```



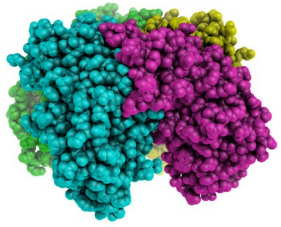
Broadcasting

```
> a = 4.  
> b = np.array([1,2,3])  
> c = np.array([[1,2,3], [4,5,6]])  
> b+a, c+a  
(array([ 5.,  6.,  7.]), array([[ 5.,  6.,  7.],  
                                [ 8.,  9., 10.])))  
> b+c  
array([[2, 4, 6],  
       [5, 7, 9]])  
> c.dot(b)  
> b.dot(c)  
ValueError: objects are not aligned  
> b[1:]*c  
ValueError: operands could not be broadcast together with shapes (2) (2,3)  
> b[1:]*c.T  
• Use matrix if you want more algebra-like behaviour
```



Broadcasting

- *The size of the trailing axes for both arrays in an operation must either be the same size or one of them must be one.*
- 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
 - they are equal, or
 - one of them is 1

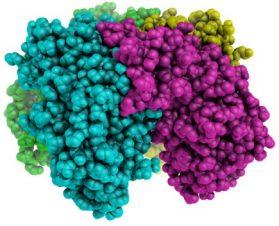


array functions and methods

- Array reduction and logical operations:

```
> a=np.arange(5)
> np.all(a>3)
False
> np.any(a>3)
True
> a > 3
array([False, False, False, False,
       True], dtype=bool)
> (a > 3) & (a < 5)
array([False, False, False, False,
       True], dtype=bool)
```

- Some details of memory use:
- `a.iscontiguous()`
- Useful when interfacing with fortran / C:
- `a.is_c_array()`
`a.is_f_array()`



Loading and saving data

- Pickle is the usual way to save and restore data in Python
- We often have data file in text format:

```
#Dist Energy
```

```
1.0 34.
```

```
1.2 38.
```

```
2.4 42.
```

- ```
> f=np.loadtxt("energies.dat")
```

```
> f
```

```
array([[1. , 34.],
 [1.2, 38.],
 [2.4, 42.]])
```

- Save single arrays with:  

```
> np.save('result_y', y)
```
- Save in text mode with:  

```
> np.savetxt('result_y', y)
```
- and multiple arrays with:  

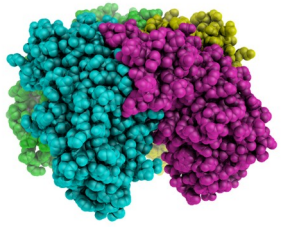
```
> np.savez('results', x, y)
```
- Recover them with load:  

```
> y=np.load('results_y.npy')
```

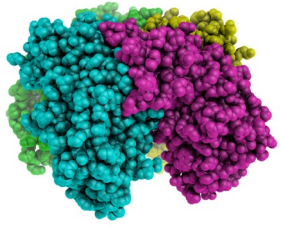
  

```
> npz=np.load('results.npz')
```



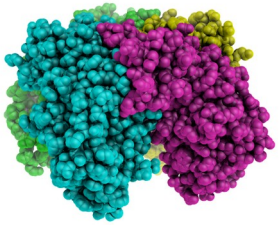


# pandas and xarray

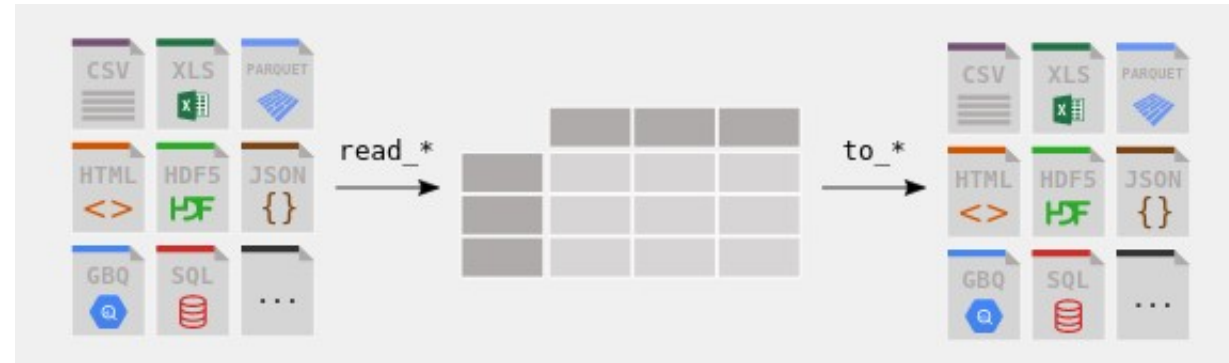


# Beyond numpy

- Pandas:
  - Dataframes with named columns and rows
  - Similar to a spreadsheet
  - Great for data analysis
  - <https://pandas.pydata.org/>
- xarray
  - labeled multidimensional arrays: n-dimensional pandas-like dataframes
  - <http://xarray.pydata.org/en/stable/>



# pandas



- import pandas as pd
- Creation of DataFrames:
  - From files (many formats):
    - read\_table, read\_csv, read\_...
  - From python dictionaries, arrays:
    - pd.DataFrame(dict)
  - From other dataframes:
    - df1.append(df2), pd.concat([df1, df2]), pd.merge

| df1 |    |    |    | df4 |   |    |    | Result |   |     |     |     |     |     |     |     |
|-----|----|----|----|-----|---|----|----|--------|---|-----|-----|-----|-----|-----|-----|-----|
|     |    |    |    |     |   |    |    |        |   |     |     |     |     |     |     |     |
|     |    |    |    |     |   |    |    |        | A | B   | C   | D   | B   | D   | F   |     |
|     | A  | B  | C  | D   |   | B  | D  | F      | 0 | A0  | B0  | C0  | D0  | NaN | NaN | NaN |
| 0   | A0 | B0 | C0 | D0  | 2 | B2 | D2 | F2     | 1 | A1  | B1  | C1  | D1  | NaN | NaN | NaN |
| 1   | A1 | B1 | C1 | D1  | 3 | B3 | D3 | F3     | 2 | A2  | B2  | C2  | D2  | B2  | D2  | F2  |
| 2   | A2 | B2 | C2 | D2  | 6 | B6 | D6 | F6     | 3 | A3  | B3  | C3  | D3  | B3  | D3  | F3  |
| 3   | A3 | B3 | C3 | D3  | 7 | B7 | D7 | F7     | 6 | NaN | NaN | NaN | NaN | B6  | D6  | F6  |
|     |    |    |    |     |   |    |    |        | 7 | NaN | NaN | NaN | NaN | B7  | D7  | F7  |

- Accessing elements:
  - df.loc[], df.iloc[]
  -

- Reshaping:
  - pd.pivot, pd.stack, pd.melt

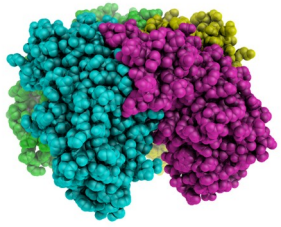
|   | foo | bar | baz | zoo |
|---|-----|-----|-----|-----|
| 0 | one | A   | 1   | x   |
| 1 | one | B   | 2   | y   |
| 2 | one | C   | 3   | z   |
| 3 | two | A   | 4   | q   |
| 4 | two | B   | 5   | w   |
| 5 | two | C   | 6   | t   |



| bar | A | B | C |
|-----|---|---|---|
| foo |   |   |   |
| one | 1 | 2 | 3 |
| two | 4 | 5 | 6 |

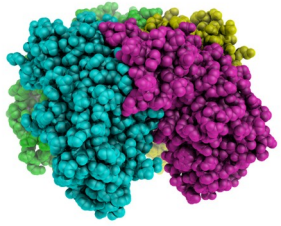
```
df.pivot(index='foo',
 columns='bar',
 values='baz')
```

- See [https://pandas.pydata.org/docs/getting\\_started/index.html](https://pandas.pydata.org/docs/getting_started/index.html)



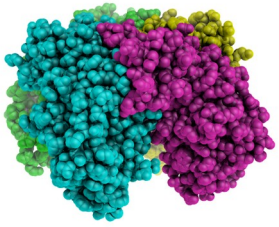
# pandas

- Compare:
  - `pd.DataFrame(np.random.random(size=(3,4)))`
  - `pd.DataFrame(np.random.random(size=(3,4)),  
                  columns=['A', 'B', 'C', 'D'])`
  - `pd.DataFrame(np.random.random(size=(3,4)),  
                  columns=['A', 'B', 'C', 'D'],  
                  index=[300., 320., 350.])`
  - `df= pd.DataFrame(np.random.random(size=(3,4)),  
                  columns=['A', 'B', 'C', 'D'],  
                  index=[300., 320., 350.])`  
`df.index.rename('Temperature', inplace=True)`



# seaborn: plotting dataframes

- `import pandas as pd`  
`import seaborn as sns`
- Seaborn is a library for making statistical graphics in Python.
  - <https://seaborn.pydata.org/introduction.html>
- `penguins = sns.load_dataset("penguins")`  
`# Show the joint distribution using kernel density estimation`  
`g = sns.jointplot(`  
    `data=penguins,`  
    `x="bill_length_mm", y="bill_depth_mm", hue="species",`  
    `kind="kde")`



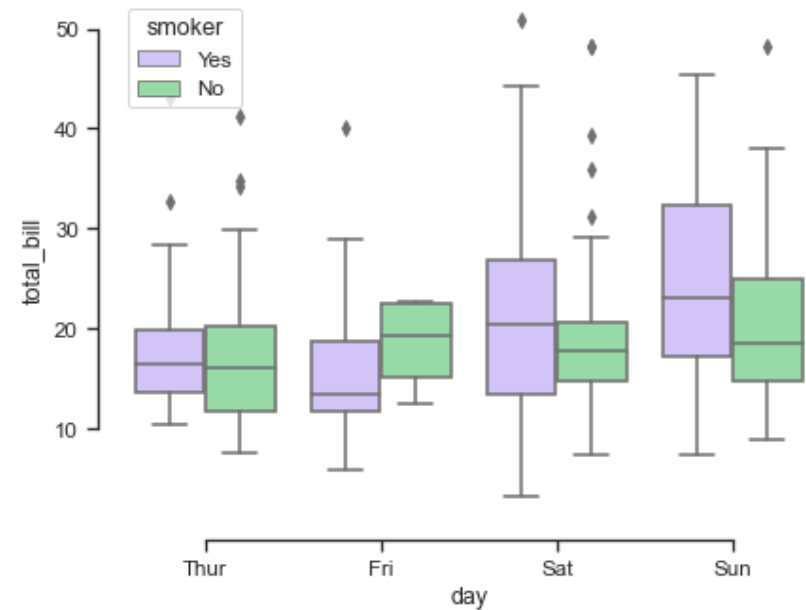
# seaborn

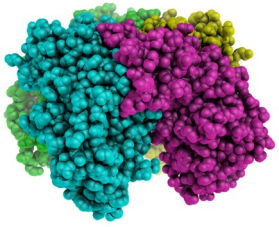
```
sns.set_theme(style="ticks",
palette="pastel")

Load the example tips dataset
tips = sns.load_dataset("tips")

Draw a nested boxplot to show bills by day and time
sns.boxplot(x="day", y="total_bill",
 hue="smoker", palette=["m", "g"],
 data=tips)

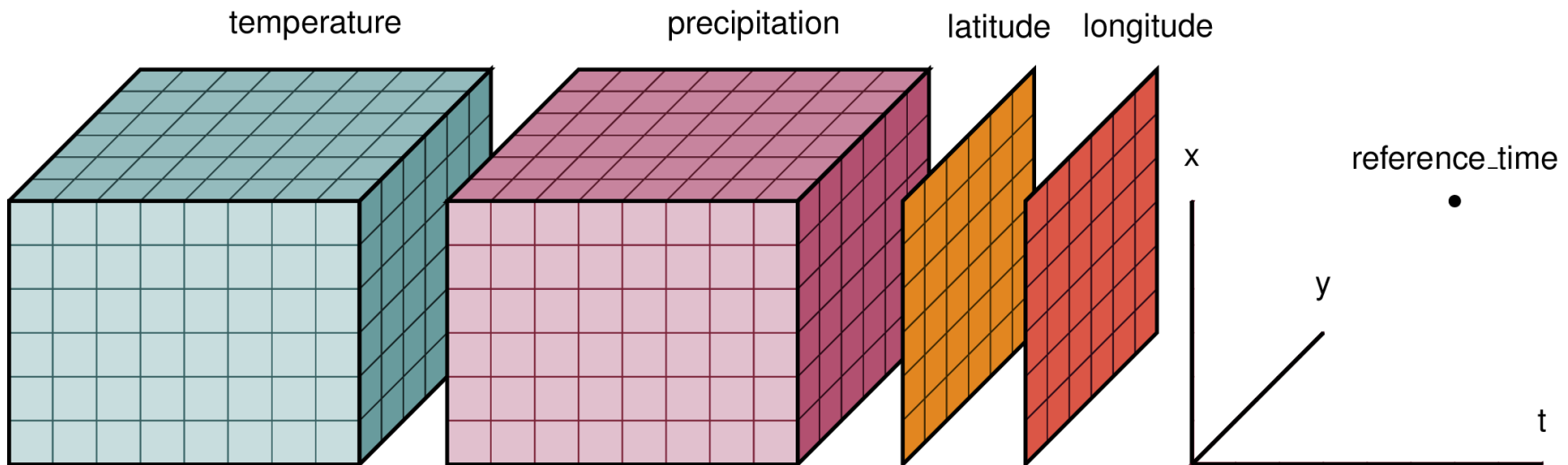
sns.despine(offset=10, trim=True)
```

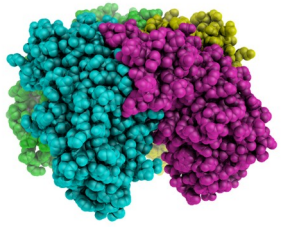




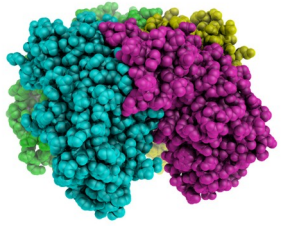
# xarray

- `import xarray as xr`
- Multi-dimensional labelled arrays (N-D pandas dataframes)
- See <http://xarray.pydata.org/en/stable/>
- ```
ds = xr.Dataset({ "temperature": ([ "x", "y", "time"], temp),  
                  "precipitation": ([ "x", "y", "time"], precip), },  
               coords={"lon": ([ "x", "y"], lon), "lat": ([ "x", "y"], lat),  
                      "time": pd.date_range("2014-09-06", periods=3),  
                      "reference_time": pd.Timestamp("2014-09-05")})
```





matplotlib

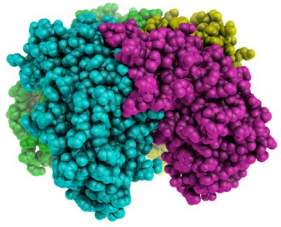


Matplotlib

- A module for plotting 2D and 3D data
- Combines well with numpy
- Starts with

```
import matplotlib.pyplot as plt  
%matplotlib inline
```

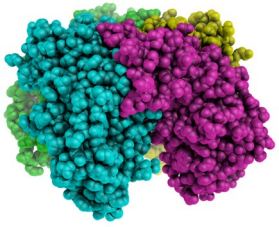
```
import pylab or similar is deprecated.
```



Matplotlib

Simplest plots:

```
> plt.plot([1,2,3], [1,4,9])  
> plt.plot(x, sin(x), '--') #where x is a numpy array  
> plt.figure() # creates new figure  
> plt.clf() # Clears current figure  
> plt.matshow(m) # m is a 2D array  
> plt.imshow(m) # m is a 2D array. Similar to matshow.  
> d = np.loadtxt('data.txt')  
> plt.plot(d[:,0], d[:,1], 's') #just slightly longer than  
gnuplot
```



Matplotlib

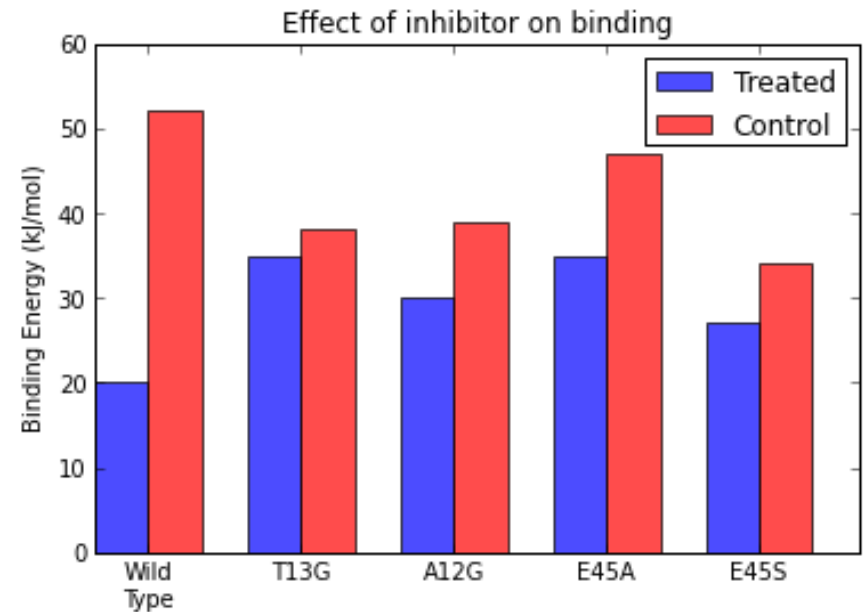
Totally reproducible
figures

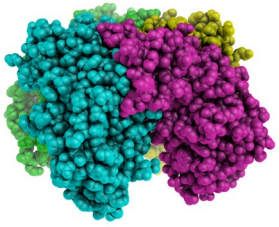
```

N = 5
treated = (20, 35, 30, 35, 27)
control = (52, 38, 39, 47, 34)
ind = np.arange(N) # the x locations for the groups
width = 0.35      # the width of the bars

fig, ax = plt.subplots()
rects1 = ax.bar(ind, treated, width, color='b', alpha=0.7, label='Treated')
rects2 = ax.bar(ind+width, control, width, color='r', alpha=0.7, label = 'Control')

# add some
ax.set_ylabel('Binding Energy (kJ/mol)')
ax.set_title('Effect of inhibitor on binding')
ax.set_xticks(ind+width)
ax.set_xticklabels( ('Wild\nType', 'T13G', 'A12G', 'E45A', 'E45S') )
ax.legend()
  
```

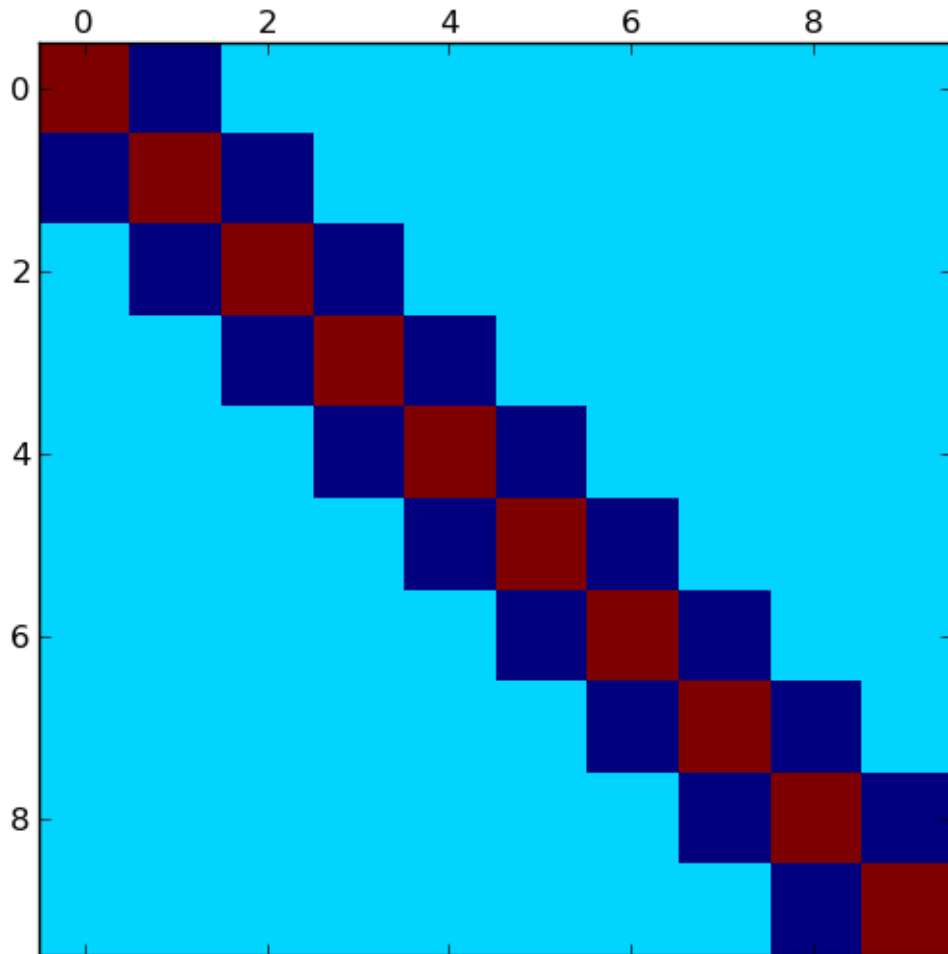




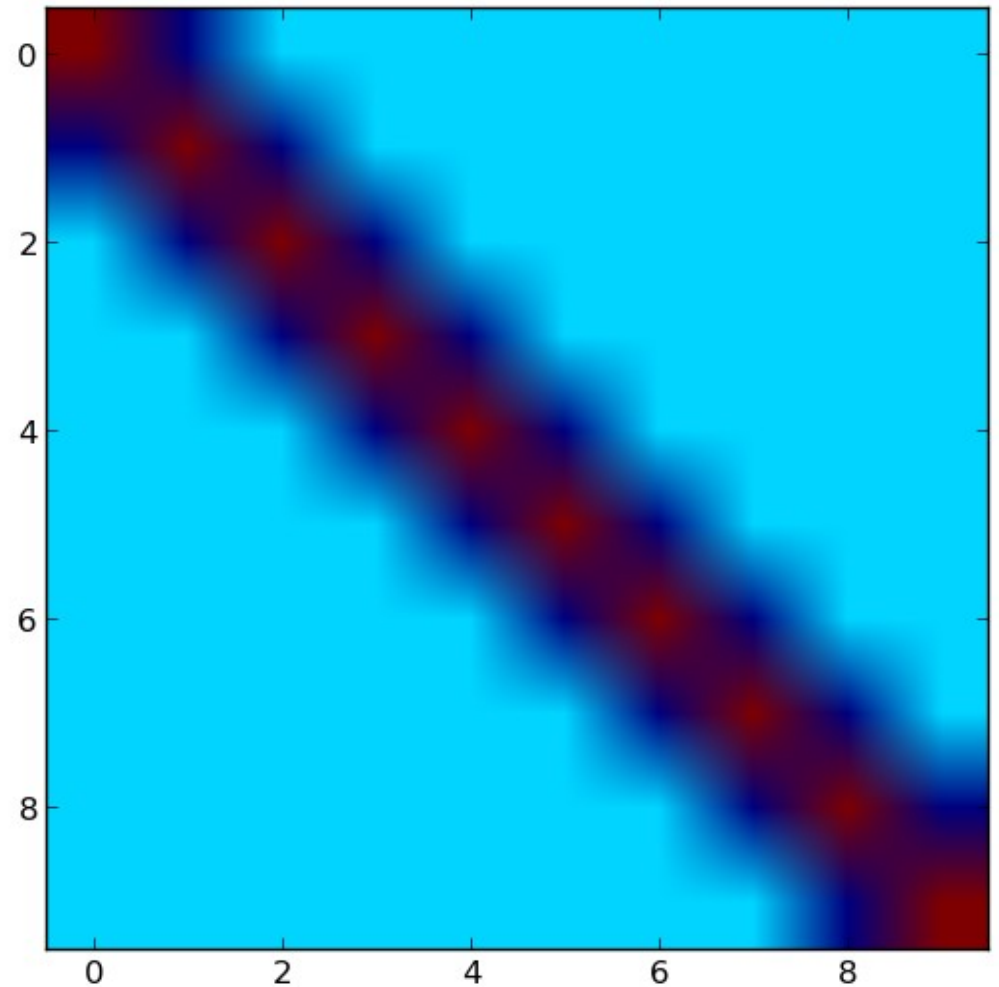
Plotting matrices

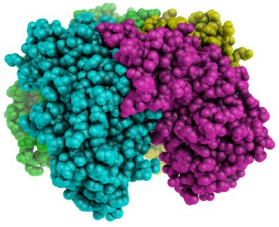
```
m=np.diag(2*np.ones(10))+np.diag(-1*np.ones(9),1)+np.diag(-1*np.ones(9), -1)
```

matshow(m)

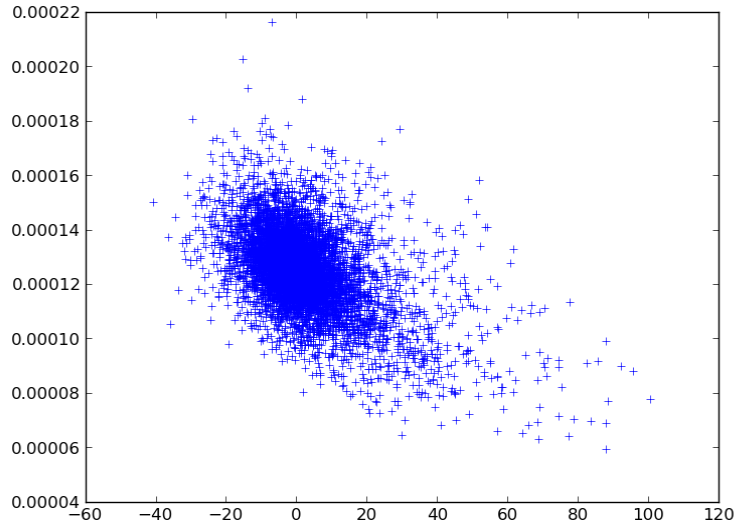


imshow(m)

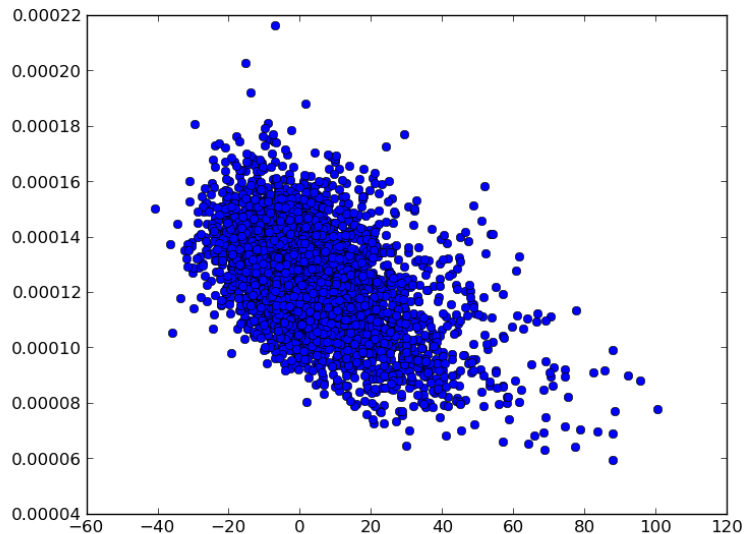




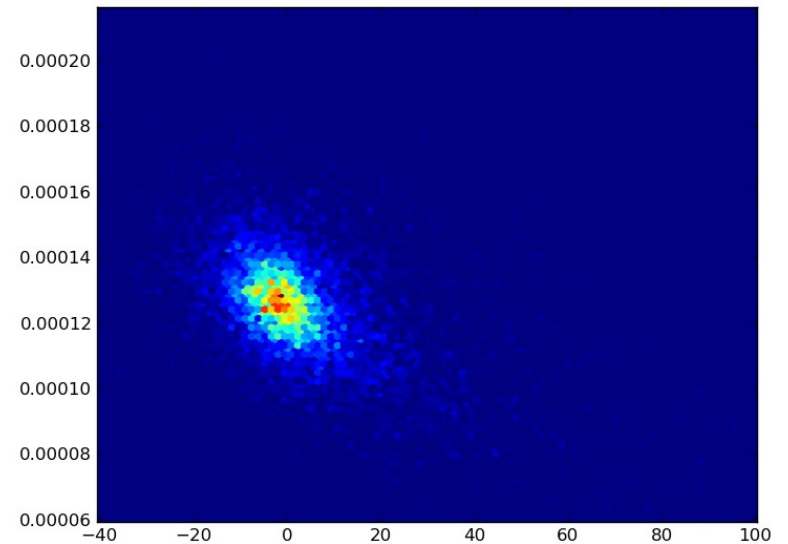
Plotting lots of points:hexbin



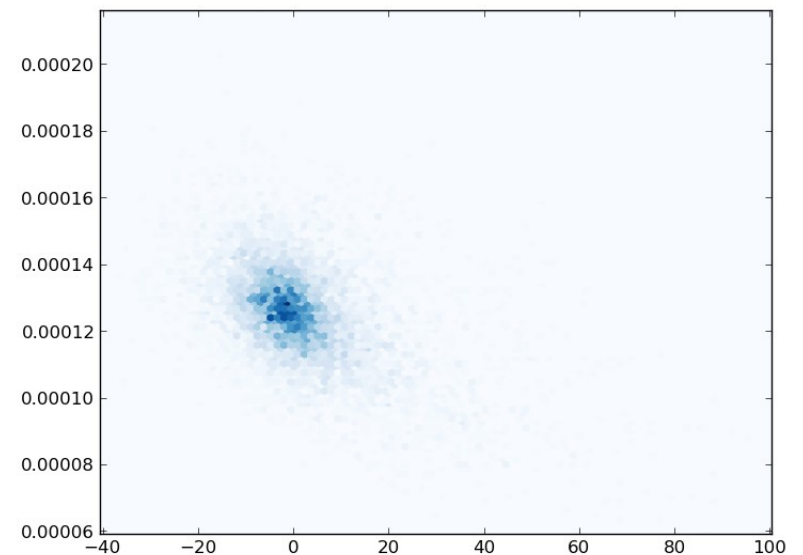
`plt.plot(x, y, '+')`



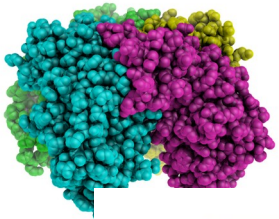
`plt.plot(x, y, 'o')`



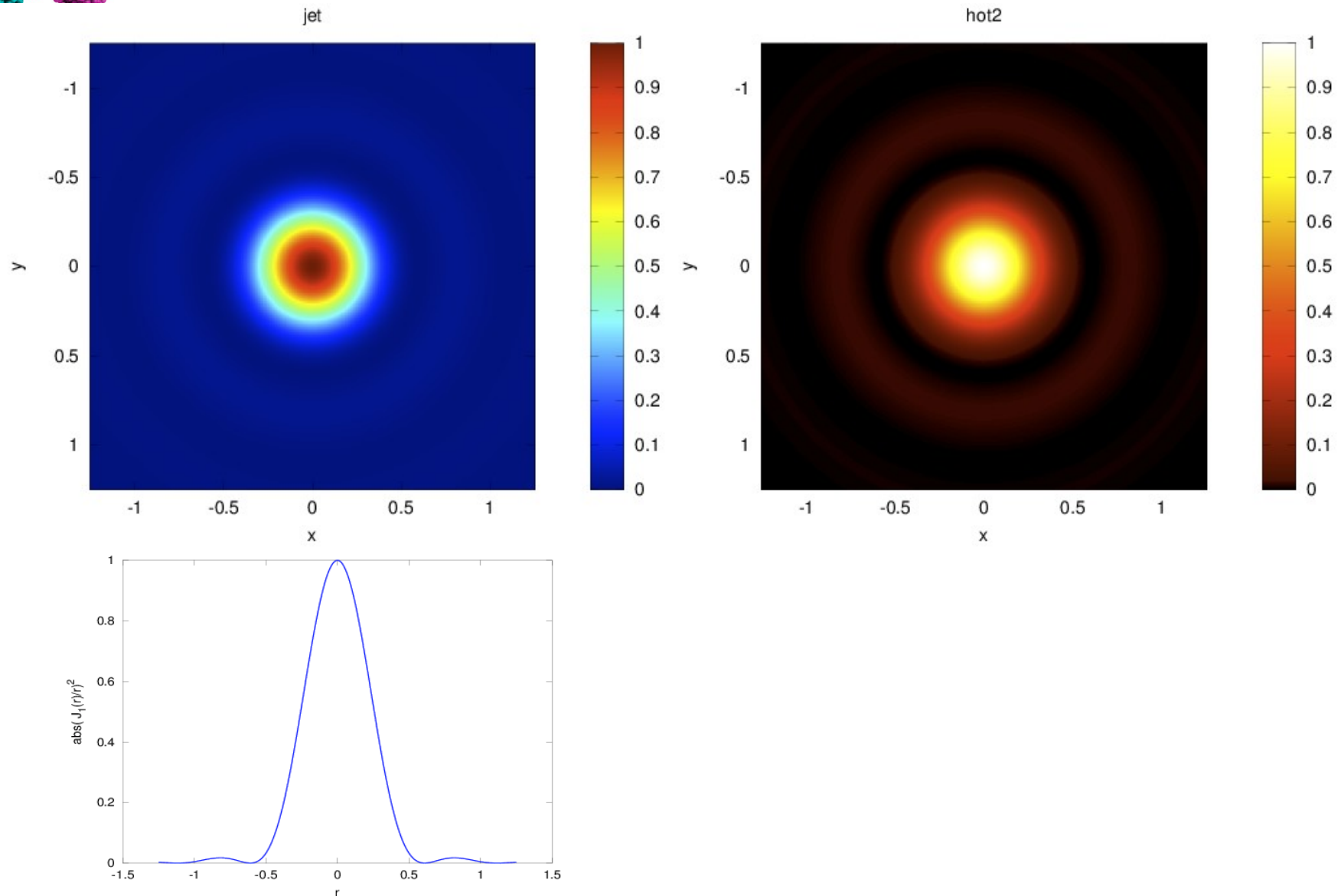
`plt.hexbin(x, y)`



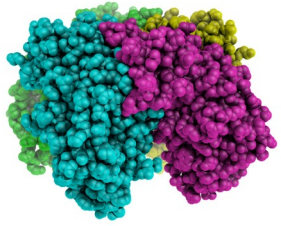
`plt.hexbin(x, y, cmap=pylab.cm.Blues)`



Jet is not a good colormap

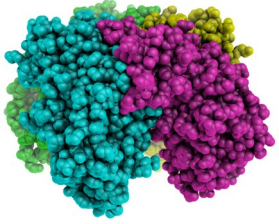


<http://cresspahl.blogspot.com.es/2012/03/expanded-control-of-octaves-colormap.html>
<https://jakevdp.github.io/blog/2014/10/16/how-bad-is-your-colormap/>



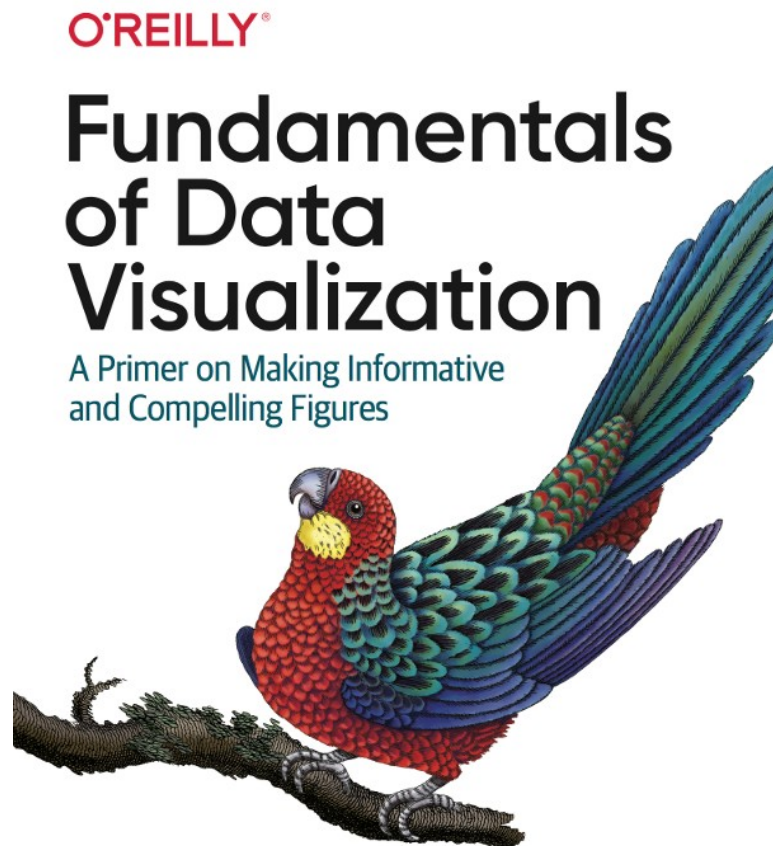
Matplotlib

- Do Lecture-4-Matplotlib.ipynb from <http://jrjohansson.github.io/>
 - Other interesting material there...
- Check matplotlib gallery
 - <http://matplotlib.org/gallery.html>
- Quick reference of symbols and colours:
 - <http://www.loria.fr/~rougier/teaching/matplotlib/#quick-references>
(part of a larger tutorial)
- Some more tricks and examples:
 - <http://wiki.scipy.org/Cookbook/Matplotlib>

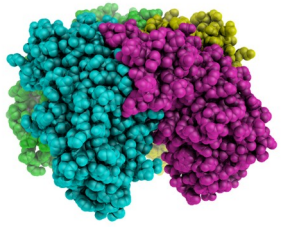


Making nice plots

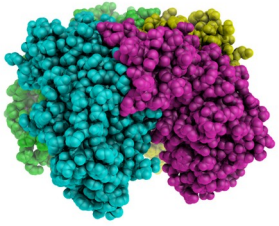
- <https://clauswilke.com/dataviz/>



Claus O. Wilke



Functions and modules



Exceptions and errors

Although the language is interpreted there are some syntax errors that prevent execution:

```
def safe_divide_1(x, y)
```

```
File"/home/ramon/python/prova.py",  
line 1
```

```
def safe_divide_1(x, y)
```

```
^
```

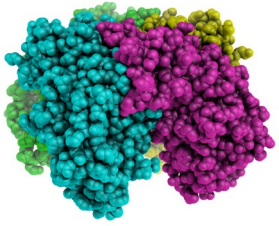
```
SyntaxError: invalid syntax
```

Exceptions leave a trace easy to follow.

Easy debugging with

```
%pdb
```

```
%debug
```



Functions

defined by def and a colon:

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

Remember indentation!

Automatic (and recommended)
documentation:

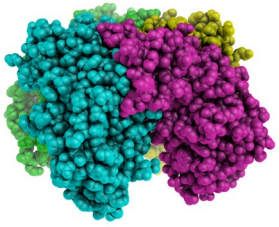
```
def add(x,y):  
    """ Returns the  
        sum of 2 numbers """  
    return x+y
```

Functions can be seen as both
fortran procedures and
functions but...

Arguments are passed by
reference

there is access to global variables:

```
> def x_val(): print(x)  
> x=60  
> x_val()  
60
```



Functions II

Function variables are local :

```
> def x_val():
```

```
...   x=40
```

```
...   print(x)
```

```
> x=60
```

```
> x_val()
```

```
40
```

```
> x
```

```
60
```

to assign variables, use return

```
def x_val():
```

```
...   x=40
```

```
...   print(x)
```

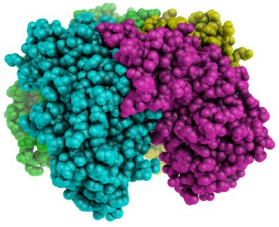
```
...   return x
```

```
> x = xval()
```

```
40
```

```
> x
```

```
40
```



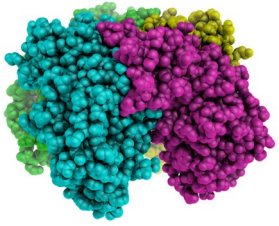
Functions III

Mutable objects are passed by reference:

```
> def square_0(lst):  
...   lst[0]*=lst[0]  
  
> a=[3,2,1]  
  
> square_0(a)  
  
> a  
[9,2,1]
```

Copy variables that need to be preserved:

```
> a_copy=a[:]  
> square_0(a)  
> import copy  
> import copy  
> a_copy=copy.deepcopy(a)
```



Functions IV

Functions can have default arguments :

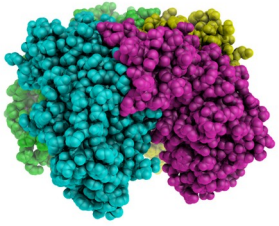
```
> def submit(job, priority=10, nprocs=1):  
...     pass  
> submit('job1.sh')
```

Function arguments do not have explicit types.

```
> add('Python ', 'summerschool')  
Python summerschool
```

Functions can be recursive

```
def fact(n):  
    if n == 1:  
        return 1  
    else:  
        return n * fact(n-1)
```



Argument unpacking

Starred arguments are tuples that collect positional arguments :

```
> def prod(*args): ...
```

```
> prod(2,3,4)
```

```
> x = (4, 5, 6)
```

```
> prod(*x)
```

In prod, args=(2,3,4)

Keyword arguments can be passed as a dictionary:

```
> options = dict(paper='A4', color = True)
```

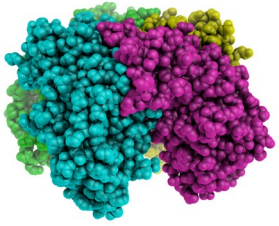
```
print_setup(options)
```

Unpacking can be a convenient way to print a list:

```
> vals = [1,2,3,4,5]
```

```
> print((4*'{:03d} ').format(*vals))
```

```
001 002 003 004
```

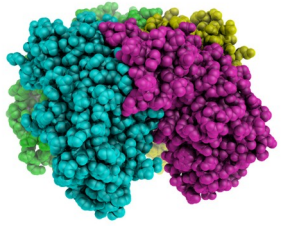


Lists or iterators?

- Lists are iterable objects
- Iterators generate objects on-the-fly
- Iterators can be created with a generator function
 - Uses `yield` statement
- Relevant for efficiency

```
def rang_llista(n):  
    result = []  
    i = 0  
    while i < n:  
        result.append(i)  
        i += 1  
    return result
```

```
def rang_gen(n):  
    i = 0  
    while i < n:  
        yield i  
        i += 1
```

Modules

- Modules allow packing libraries or extensions
- There are built-in and external modules
- When imported modules are executed
- Modules can be written in C or Fortran!

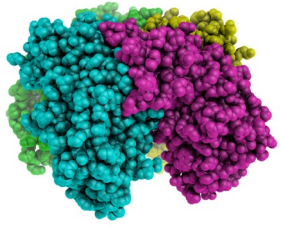
```
> import math
```

```
> m = math
```

```
> import math as m
```

```
> from math import cos, sin
```

```
> from math import * #dangerous. All into the same namespace
```

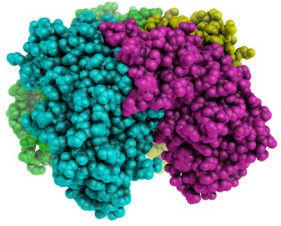


Modules

- Python checks if a module is already loaded.
 - The interpreter does not reload a module already imported
 - This can cause unexpected behaviour interactively
- Ipython has a more versatile module loading

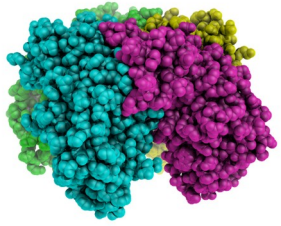
```
%load_ext autoreload
```

```
autoreload 2 #Will reload a module if it changes
```



Some useful modules

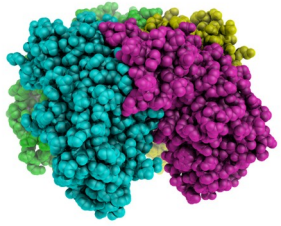
- `sys` — System-specific parameters and functions
- `os` — Miscellaneous operating system interfaces
- `os.path` — Common pathname manipulations
- `glob` — Unix style pathname pattern expansion
- `re` — regular expressions
- `copy` — Shallow and deep copy operations
- `argparse` — Parser for command-line options, arguments and sub-commands
- `subprocess` — Subprocess management
- `inspect` — Inspect live objects



Some useful modules

```
if len(sys.argv)!=3:
    print('Error: Use two arguments.')
    sys.exit()

method = sys.argv[1]
filelist = glob.glob('/home/ramon/*')
for fileName in filelist:
    if os.path.isfile(fileName): print(fileName)
```

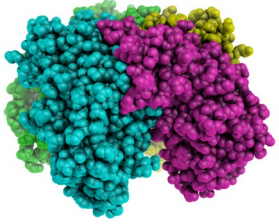


Modules: too many...

From the python documentation:

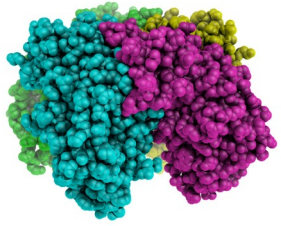
It is also possible to use a list as a queue, where the first element added is the first element retrieved (“first-in, first-out”); however, lists are not efficient for this purpose. While appends and pops from the end of list are fast, doing inserts or pops from the beginning of a list is slow (because all of the other elements have to be shifted by one).

To implement a queue, use `collections.deque` which was designed to have fast appends and pops from both ends.



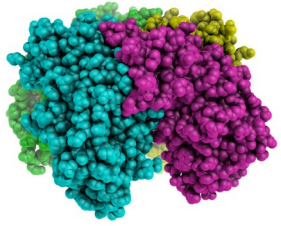
Modules: too many...

```
>>> import math
>>> import cmath
>>> import numpy.lib.scimath as scimath
>>> math.sqrt(4)
2.0
>>> math.sqrt(-4)
Traceback (most recent call last):
  File "<stdin>", line 1, in <module>
ValueError: math domain error
>>> cmath.sqrt(4)
(2+0j)
>>> cmath.sqrt(-4)
2j
>>> scimath.sqrt(4)
2.0
>>> scimath.sqrt(-4)
2j
```



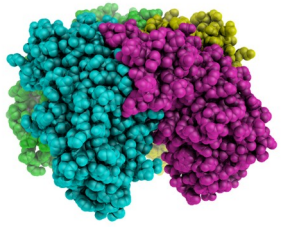
Working with your modules

- Import reads from local directory and from the directories in `sys.path` (import `sys` first)
- Put your modules in a directory and add it to the environment variable `$PYTHONPATH`.
- Python will add the directories in `$PYTHONPATH` to `sys.path`
- Document your modules and the functions therein.
- Use `if __name__ == '__main__':` to execute code only if Python is running the module, and not if it is imported.
 - <http://stackoverflow.com/questions/419163/what-does-if-name-main-do>

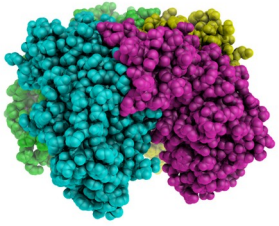


Installing external Modules

- Use conda distribution. Then ``conda install module``
- Many come as part of the linux distributions (usually older versions)
 - ipython, numpy, biopython...
- For modules in the PyPI repository (most of them)
<https://pypi.python.org/pypi>
 - Use `pip-3` or `pip3`
- Use:
 - `$ python setup.py build`
 - `$ (sudo?) python setup.py install`



Scipy



Linear algebra

- Support for LAPACK, BLAS and ATLAS
 - Can make Scipy compilation more involved

```
> A=matrix(random.rand(5,5))
```

```
> A.I
```

```
> linalg.det(A)
```

```
> linalg.eigvals(A)
```

```
> linalg.eig(A)
```

```
> linalg.svd(A)
```

```
> linalg.cholesky(A)
```

- Solving linear systems:

- $\mathbf{A} \cdot \mathbf{x} = \mathbf{b}$

```
>
```

```
b=matrix(random.rand(5)).reshape((5,1))
```

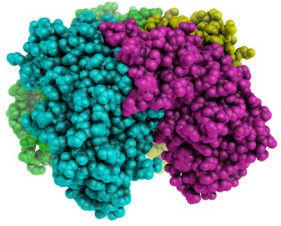
```
> linalg.solve(A,b)
```

- LAPACK, BLAS wrappers

```
> from scipy.lib import lapack
```

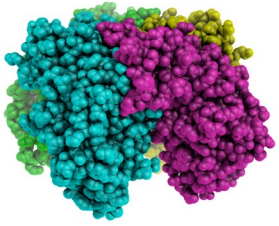
```
> from scipy.lib import blas
```

```
blas.fblas.sdot?
```



Linear regression

- There are different ways to perform linear regression (still surprised?)
- See the notebook.



Optimization

- There are different optimization methods:

```
> import scipy.optimize as so
```

- Some only need the function value:

```
> fmin, fmin_powell
```

- Some need the gradient or the hessian:

```
> fmin_cg, fmin_bfgs, fmin_ncg
```

- Some look for global minima:

```
> anneal
```

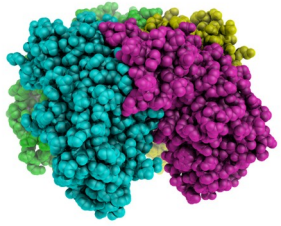
- Remember:

```
> scipy.info('optimize')
```

- Pedagogical documentation:

- <http://docs.scipy.org/doc/scipy/reference/tutorial/optimize.html>

- <http://docs.scipy.org/doc/scipy/reference/optimize.html>

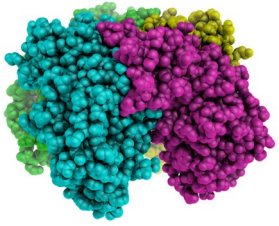


f2py

- Many things are fast with Numpy
- Iterative algorithms over **array values** are slow
- You can import Fortran functions and subroutines with f2py
- You could also call external fortran programs with

```
> subprocess.call(<program>, shell=True)
```

 - but data exchange has to be through files (slower)
- f2py finds your fortran compiler. Works with gfortran, ifort,...
- f2py creates a module you can import in python
- As simple as:
 - `$ f2py -c <file> -m <module>`
 - Tip: first compile it to check it works



f2py II

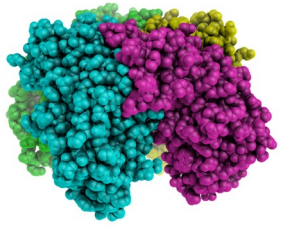
```
module funcs
implicit none
contains
function f1(x,y)
  real,intent(in):: x,y
  real:: f1
  f1=x+y**2
end function f1

function f2(x,y)
  real,intent(in):: x,y
  real, dimension(3):: f2
  f2(1)=x+y**2
  f2(2)=sin(x*y)
  f2(3)=2*x-y
end function f2
end module
```

```
$ f2py -c test.f90 -m test
```

- go to ipython:

```
> import test
> test.funcs.f1(1,2)
5.0
> test.funcs.f2(1,2)
array([ 5., 0.90929741, 0.],
      dtype=float32)
```



f2py III

Using ipython magicfunctions:

```
sudo pip3 install -U fortran-magic
```

Useful for performing long array operations

```
In [5]: %load_ext fortranmagic
```

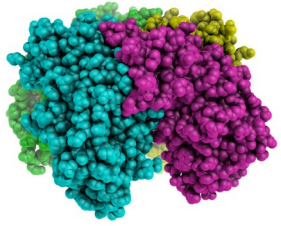
```
In [6]: %%fortran
        subroutine f1(x, y, z)
            real, intent(in) :: x,y
            real, intent(out) :: z

            z = sin(x+y)

        end subroutine f1
```

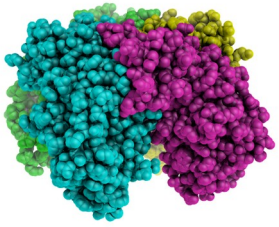
```
In [7]: f1(1.0, 2.1415)
```

```
Out[7]: 9.26574066397734e-05
```



Big data, big memory

- Numpy arrays are meant to live in memory
- If that is not possible:
 - Use `op=` operations (they use half the memory):
 - `p *= alpha` is better than `p = p * alpha`
 - Use `scipy.sparse` matrices
 - Use `PyTable` to store (compressed) matrices on disk
 - Modify your algorithm to work with submatrices



Sympy: Symbolic math

- Symbolic algebra
- Analytic solution of equations
- Integration, derivation
- Polynomials
- Limis
-

Alternate forms:

`(cos(x + y)).expand(trig=True)`

$$-\sin(x)\sin(y) + \cos(x)\cos(y)$$

`trigsimp(cos(x + y))`

$$\cos(x + y)$$

`(cos(x + y)).rewrite(csc, sin, sec, cos, cot, tan)`

$$\frac{-\tan^2\left(\frac{x}{2} + \frac{y}{2}\right) + 1}{\tan^2\left(\frac{x}{2} + \frac{y}{2}\right) + 1}$$

`(cos(x + y)).rewrite(sin, exp, cos, exp, tan, exp)`

$$\frac{1}{2}e^{i(-x-y)} + \frac{1}{2}e^{i(x+y)}$$

<http://sympy.org/en/index.htm>

|

```
>>> integ = Integral(sin(x**2), x)
```

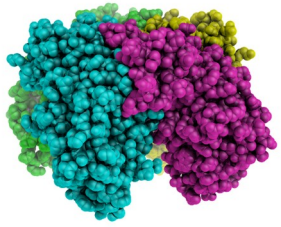
```
>>> integ
```

$$\int \sin(x^2) dx$$

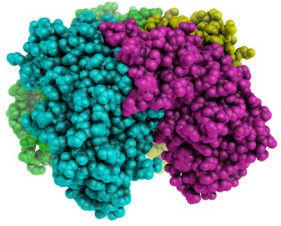
```
>>> integ.doit()
```

$$3 \cdot \sqrt{2} \cdot \sqrt{\pi} \cdot \text{fresnels}\left(\frac{\sqrt{2} \cdot x}{\sqrt{\pi}}\right) \cdot \Gamma(3/4)$$

$$8 \cdot \Gamma(7/4)$$



Classes



Classes and objects

Most simple examples are quite stupid...

Everything is an object with python.

Objects have methods

`sort` is a method of list objects

They have to be called with `()`

Objects have attributes

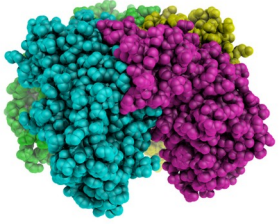
`c.real` is an attribute

Example: you have a set of sensors

Sensors have position

Sensors have a value

Sensors can be reset



Classes and objects II

Sensor is an object

Each object has methods (and attributes)

```
sensor[i].position
```

```
sensor[i].value
```

```
sensor[i].reset(0.0)
```

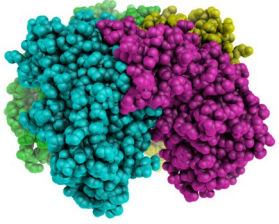
Fortran / C:

Generate arrays for each property:

```
position[i]
```

```
value[i]
```

```
subroutine reset(i, val)
```



Classes and objects III

Better Fortran / C:

use struct (in C)

use type (in Fortran)

type sensor

 real, dimension(3) ::
position

 real :: value

end type sensor

sensor[i]%position

sensor[i]%value

reset ?

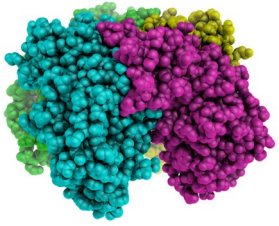
But now you also have clocks:

clocks have values

clocks can be reset

What does value refer to?

What is the argument of reset?



Objects and inheritance

New classes are defined as:

```
class Point:
```

```
    def __init__(self,x,y):
```

```
        self.x=x
```

```
        self.y=y
```

Classes allow the definition of elaborate personal objects

```
class Num_List(list):
```

```
    def __init__(self, lst=None)
```

```
        list.__init__(self,lst)
```

```
    def square(self):
```

```
        for i in range(len(self)):
```

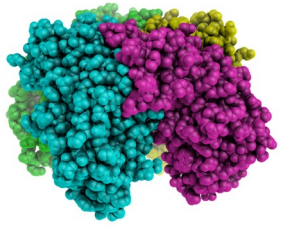
```
            self[i]=self[i]**2
```

```
> a=num_list([1,2,3])
```

```
> a.square()
```

```
> a
```

```
[1, 4, 9]
```



A periodic table of objects

Each chemical element can be defined as a object.

Attributes such as:

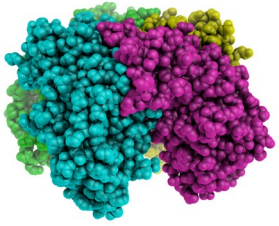
- Name
- atomic symbol
- atomic weight
- Exercice: generate element objects and iterate over them to print:

The atomic number of hydrogen is 1

The atomic number of lithium is 3

The atomic number of helium is 2

- <https://ramoncrehuet.wordpress.com/2014/11/12/a-periodic-table-of-objects/>



A practical application

Line search in one dimension:

$f(x)$ becomes $f(\mathbf{x}_0 + s \mathbf{d})$

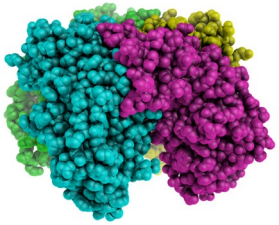
Implicitly depends on \mathbf{x}_0 and \mathbf{d} .

How can we use our line_search algorithm?

```
def f1(x):  
    return x[0]**2+2*x[1]**2+4*\  
           x[0]*x[1]-3*x[1]+2*x[1]
```

```
x0 = np.asarray([3,-1.])  
r=cg(f1, x0)  
print(r)
```

```
import numpy as np  
import scipy.optimize as so  
def cg(func, x0):  
    "Conjugate Gradient"  
    class Scalar:  
        'A class to hold scalar functions'  
        def __init__(self, func, x0, d):  
            self.func = func  
            self.x0 = x0  
            self.d = d  
        def value(self,s):  
            'function evaluation'  
            return self.func(x0+s*d)  
    d = # from conjugate gradient algorithm  
    f=Scalar(func, x0, d)  
    result = so.minimize_scalar(f.value)  
    return result
```

A practical application

As usual, there are already solutions for this common task:

```
scipy.optimize.fmin(func, x0, args=())
```

Parameters

`func` : callable `func(x,*args)`

The objective function to be minimized.

`x0` : ndarray

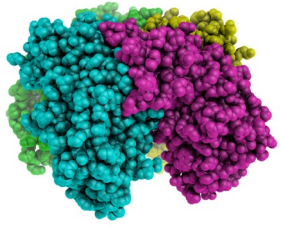
Initial guess.

`args` : tuple, optional

Extra arguments passed to `func`, i.e. ```f(x,*args)```.

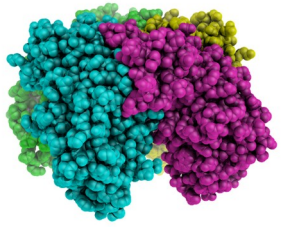
Or use (needs gradient):

```
scipy.optimize.line_search
```

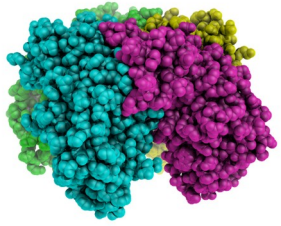


Add ons: itertools

```
> import itertools
> perms = itertools.permutations('ABC', 3)
> list(perms)
[('A', 'B', 'C'),
 ('A', 'C', 'B'),
 ('B', 'A', 'C'),
 ('B', 'C', 'A'),
 ('C', 'A', 'B'),
 ('C', 'B', 'A')]
> list(itertools.combinations('ABC',2))
[('A', 'B'), ('A', 'C'), ('B', 'C')]
```



Resources



Resources

On-line Official documentation (contains Tutorial in PDF or HTML):

<http://www.python.org/doc>

General introductory books (also in paper):

<http://diveintopython.org/> (This one is simpler!)

<http://www.greenteapress.com/thinkpython/thinkpython.html>

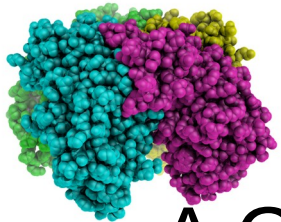
Comparison of codes in different languages:

<http://rosetacode.org>

<http://www.codecodex.com>

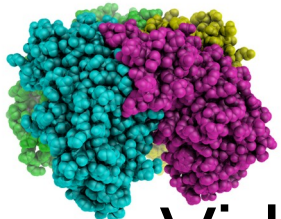
Python package index: where to find modules

<http://pypi.python.org/pypi>



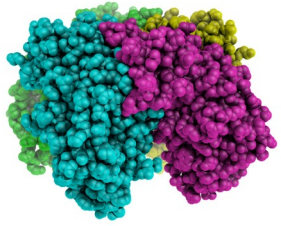
Resources

- A Crash Course in Python for Scientists (with applications in Quantum chemistry)
 - <http://nbviewer.ipython.org/5920182>
 - Written in an ipython notebook
- Python Scientific Lectures
 - <http://scipy-lectures.org/>
- Python flow with Pythontutor
 - <http://www.pythontutor.com>



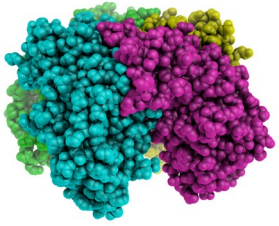
Resources

- Videos from the Scipy conference:
 - https://www.youtube.com/watch?v=V0D2mhVt7NE&t=0s&list=PLYx7XA2nY5Gd-tNhm79CNMe_qvi35PgUR&index=15
 - https://www.youtube.com/watch?v=Gzun8PpyBCo&t=0s&list=PLYx7XA2nY5Gd-tNhm79CNMe_qvi35PgUR&index=93
 -



Resources: Books

- Langtangen, *A Premier on Scientific Programming with Python*.
 - Good for learning scientific programming. Starts from scratch. Unfortunately does not use ipython nor numpy. Python 2.
- Langtangen, *Python Scripting for Computational Science*.
 - Good book for programmers. Advanced level. Explains how to interface with C and Fortran and how to optimize code. Python 2.
- Rossant, C, *Learning Ipython for Interactive Computing and Data Visualization*.
 - Basic level. Covers several subjects, including matplotlib and parallelism.



Resources: Teaching

- On teaching programming with Python 3

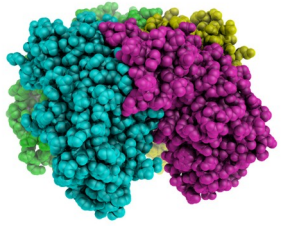
<http://www.comp.leeds.ac.uk/nde/papers/teachpy3.html>

- Online Syntax Highlighting

<http://tohtml.com/python/>

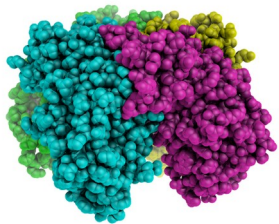
- Style Guide for Python Code:

- www.python.org/dev/peps/pep-0008/



K. Hinsen views

- NumPy has introduced incompatible changes with almost every new version over the last years
- Even the importance of NumPy in the scientific Python ecosystem, I consider its lack of stability alarming.
- What makes me hesitate to recommend not using Python is that there is no better alternative.
- <https://khinsen.wordpress.com/2014/09/12/the-state-of-numpy/>



Software in python

- Molecular visualization:
 - VMD: <http://www.ks.uiuc.edu/Research/vmd/>
 - pymol: <http://www.pymol.org/>
- Molecular Dynamics
 - Espresso <http://espressomd.org/wordpress/>
 - HOOMD-Blue <http://glotzerlab.engin.umich.edu/hoomd-blue/>
 - QM/MM with pDynamo: <http://www.pdynamo.org>
- QM calculation with
 - pyQuante: <http://pyquante.sourceforge.net/>
 - NWChem: <http://www.nwchem-sw.org/index.php/Python>
- Protein structure with pyRosetta: <http://pyrosetta.org/>
- Bioinformatics with BioPython: <http://biopython.org/>