

4. numpy

May 5, 2014

1 La biblioteca numpy: “numerical Python” (Python numérico)

```
In [1]: import numpy
```

Numpy provee *vectores* (en el sentido matemático)

Recordemos que al sumar dos listas, no hace lo que queremos que haga:

```
In [2]: l = [1, 2, 3]
        l + l
```

```
Out[2]: [1, 2, 3, 1, 2, 3]
```

```
In [3]: numpy.array # es una funcion que acepta una lista
```

```
Out[3]: <function numpy.core.multiarray.array>
```

```
In [6]: v = numpy.array(1)
```

Acepta una lista y regresa un `array` de numpy

```
In [7]: v + v
```

```
Out[7]: array([2, 4, 6])
```

Ahora sí se comporta como un vector!

```
In [9]: 3.5 * v # multiplica cada elemento
```

```
Out[9]: array([ 3.5,  7. , 10.5])
```

```
In [10]: numpy.array([2., 1.] )
```

```
Out[10]: array([ 2.,  1.])
```

```
In [11]: dot(v, v)
```

```
-----
NameError
```

```
Traceback (most recent call last)
```

```
<ipython-input-11-093f352c32e0> in <module>()
----> 1 dot(v, v)
```

```
NameError: name 'dot' is not defined
```

```

In [12]: %who

a      1      numpy      v

In [13]: numpy.dot(v, v)

Out[13]: 14

In [14]: numpy.dot

Out[14]: <function numpy.core.multiarray.dot>

In [15]: v

Out[15]: array([1, 2, 3])

In [16]: type(v)

Out[16]: numpy.ndarray

    Matrices:

In [18]: M = numpy.array([ [2., 1.], [1., 1.] ])

In [19]: M

Out[19]: array([[ 2.,  1.],
                [ 1.,  1.]])

In [20]: print(M)

[[ 2.  1.]
 [ 1.  1.]]

In [21]: M + M

Out[21]: array([[ 4.,  2.],
                [ 2.,  2.]])

In [23]: M * M # atención: componente por componente; NO es multiplicacion matricial normal

Out[23]: array([[ 4.,  1.],
                [ 1.,  1.]])

In [25]: numpy.dot(M, M)

Out[25]: array([[ 5.,  3.],
                [ 3.,  2.]])


$$(AB)_{ij} = \sum_k A_{ik} B_{kj}$$

Es una contracción de índices
Normalmente:

In [26]: import numpy as np

In [27]: np.array([4, 5, 5])

Out[27]: array([4, 5, 5])

```

No usemos `from numpy import *` – importa ~500 funciones
numpy tiene cualquier operación sobre vectores y matrices (y su equivalente en más dimensiones – tensores de más alto rango)
Unos ejemplos:

```
In [28]: np.zeros(5)
```

```
Out[28]: array([ 0.,  0.,  0.,  0.,  0.])
```

```
In [29]: np.zeros( (10, 10) )  # (10, 10) es un par ordenado (tupla)
```

```
Out[29]: array([[ 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.,  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.,  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 [30]: (10, 10)
```

```
Out[30]: (10, 10)
```

```
In [31]: np.ones(10)
```

```
Out[31]: array([ 1.,  1.,  1.,  1.,  1.,  1.,  1.,  1.,  1.,  1.])
```

```
In [33]: 2. * np.ones(10)
```

```
Out[33]: array([ 2.,  2.,  2.,  2.,  2.,  2.,  2.,  2.,  2.,  2.])
```

```
In [34]: np.identity(5)
```

```
Out[34]: array([[ 1.,  0.,  0.,  0.,  0.],
                 [ 0.,  1.,  0.,  0.,  0.],
                 [ 0.,  0.,  1.,  0.,  0.],
                 [ 0.,  0.,  0.,  1.,  0.],
                 [ 0.,  0.,  0.,  0.,  1.]])
```

```
In [35]: np.diag([1, 2, 3])
```

```
Out[35]: array([[1, 0, 0],
                 [0, 2, 0],
                 [0, 0, 3]])
```

1.1 Cómo extraer elementos de una matriz:

```
In [40]: M = np.diag(range(5))
         M
```

```
Out[40]: array([[0, 0, 0, 0, 0],
                 [0, 1, 0, 0, 0],
                 [0, 0, 2, 0, 0],
                 [0, 0, 0, 3, 0],
                 [0, 0, 0, 0, 4]])
```

```

In [41]: M = np.array([1,2,3,4])
In [42]: M
Out[42]: array([1, 2, 3, 4])
In [43]: M.shape
Out[43]: (4,)
In [45]: M = np.array([1,2,3,4]).reshape(2, 2)
          M
Out[45]: array([[1, 2],
               [3, 4]])
In [47]: M[0] # primer renglon
Out[47]: array([1, 2])
In [48]: M[0][1]
Out[48]: 2
In [49]: M[0, 1]
Out[49]: 2
In [54]: M[0, 0:2]
Out[54]: array([1, 2])
In [55]: M[0, 0:-1]
Out[55]: array([1])
In [56]: M[0, :] # toma por defecto el inicio y fin del vector
Out[56]: array([1, 2])
In [58]: M[:, 0] # primera columna
Out[58]: array([1, 3])
In [60]: M.T # da la transpuesta
Out[60]: array([[1, 3],
               [2, 4]])
In [61]: M.T[0]
Out[61]: array([1, 3])
In [62]: M[:,0]
Out[62]: array([1, 3])
In [65]: M = np.diag(range(5))
In [66]: M
Out[66]: array([[0, 0, 0, 0, 0],
               [0, 1, 0, 0, 0],
               [0, 0, 2, 0, 0],
               [0, 0, 0, 3, 0],
               [0, 0, 0, 0, 4]])
In [68]: M[1:4, 2:4] # submatriz
Out[68]: array([[0, 0],
               [2, 0],
               [0, 3]])

```

1.2 Álgebra lineal

Hay un submódulo de `numpy` que se llama `linalg` que provee las funciones básicas de álgebra lineal:

```
In [69]: from numpy import linalg
```

```
In [70]: linalg
```

```
Out[70]: <module 'numpy.linalg' from '/usr/local/Cellar/python/2.7.5/Frameworks/Python.framework/Versions/2.7/lib/python2.7/site-packages/numpy/linalg.py'>
```

```
In [71]: M
```

```
Out[71]: array([[0, 0, 0, 0, 0],
                [0, 1, 0, 0, 0],
                [0, 0, 2, 0, 0],
                [0, 0, 0, 3, 0],
                [0, 0, 0, 0, 4]])
```

```
In [72]: linalg.inv(M)
```

```
-----
LinAlgError
```

```
Traceback (most recent call last)
```

```
<ipython-input-72-8f8888e69724> in <module>()
----> 1 linalg.inv(M)
```

```
/usr/local/Cellar/python/2.7.5/Frameworks/Python.framework/Versions/2.7/lib/python2.7/site-packages/numpy/linalg.py
443     """
444     a, wrap = _makearray(a)
--> 445     return wrap(solve(a, identity(a.shape[0], dtype=a.dtype)))
446
447
```

```
/usr/local/Cellar/python/2.7.5/Frameworks/Python.framework/Versions/2.7/lib/python2.7/site-packages/numpy/linalg.py
326     results = lapack_routine(n_eq, n_rhs, a, n_eq, pivots, b, n_eq, 0)
327     if results['info'] > 0:
--> 328         raise LinAlgError('Singular matrix')
329     if one_eq:
330         return wrap(b.ravel().astype(result_t))
```

```
LinAlgError: Singular matrix
```

```
In [73]: linalg.det(M)
```

```
Out[73]: 0.0
```

```
In [74]: M = np.array([1, 2, 3, 4])
```

```
In [77]: N = M.reshape(2, 2) # es una vista ("view") de la matriz original
        # NO es una copia; son los mismos datos
```

```
In [78]: N
```

```

Out[78]: array([[1, 2],
               [3, 4]])

In [79]: N[0, 1] = 100

In [80]: N

Out[80]: array([[ 1, 100],
               [ 3,  4]])

In [81]: M

Out[81]: array([ 1, 100,  3,  4])

In [82]: N

Out[82]: array([[ 1, 100],
               [ 3,  4]])

In [83]: N[0, 1] = 2

In [83]:

In [84]: N

Out[84]: array([[1, 2],
               [3, 4]])

In [86]: linalg.det(N)

Out[86]: -2.0000000000000004

In [87]: linalg.inv(N)

Out[87]: array([[-2. ,  1. ],
               [ 1.5, -0.5]])

In [89]: linalg.inv??

In [91]: linalg.eig(N)

Out[91]: (array([-0.37228132,  5.37228132]), array([[-0.82456484, -0.41597356],
               [ 0.56576746, -0.90937671]]))

```

Regresa una tupla (par ordenado) de dos elementos: e-vals y e-vecs

```

In [93]: lamb, v = linalg.eig(N)
         # tupla = tupla -- hace desempaque de tuplas

In [94]: lamb

Out[94]: array([-0.37228132,  5.37228132])

In [95]: v

Out[95]: array([[-0.82456484, -0.41597356],
               [ 0.56576746, -0.90937671]])

```

Ejercicio: Averigua dónde están los eigenvectores

Tarea: Método de potencias para calcular un eigenvector y un eigenvalor

```
In [6]: import numpy as np
        from numpy import linalg

In [3]: M = np.array([2., 1., 1., 1.]).reshape(2, 2)

In [4]: c = np.array([1., 1.])
```

Quiero resolver el sistema $M \cdot x = c$:

```
In [8]: x = linalg.solve(M, c)
        x
```

```
Out[8]: array([ 0.,  1.])
```

```
In [13]: linalg.solve?
```

1.3 Transformadas de Fourier

```
In [97]: from numpy import fft # fast Fourier transform
```

```
In [99]: x = np.linspace(0, 100, 1024)
```

```
In [100]: x
```

```
Out[100]: array([ 0.00000000e+00,  9.77517107e-02,  1.95503421e-01, ...,
                  9.98044966e+01,  9.99022483e+01,  1.00000000e+02])
```

```
In [101]: len(x)
```

```
Out[101]: 1024
```

```
In [102]: f = fft.fft(x)
```

```
In [103]: x2 = fft.ifft(f)
```

```
In [104]: x2
```

```
Out[104]: array([ 3.55271368e-15 -2.49800181e-15j,
                  9.77517107e-02 -2.22044605e-16j,
                  1.95503421e-01 -5.01889652e-15j, ...,
                  9.98044966e+01 +3.88578059e-15j,
                  9.99022483e+01 +1.82027949e-14j,  1.00000000e+02 -1.17267307e-14j])
```

```
In [105]: x
```

```
Out[105]: array([ 0.00000000e+00,  9.77517107e-02,  1.95503421e-01, ...,
                  9.98044966e+01,  9.99022483e+01,  1.00000000e+02])
```

```
In [107]: x2.real
```

```
Out[107]: array([ 3.55271368e-15,  9.77517107e-02,  1.95503421e-01, ...,
                  9.98044966e+01,  9.99022483e+01,  1.00000000e+02])
```

```
In [108]: x2.real - x # errores
```

```
Out[108]: array([ 3.55271368e-15,  7.02216063e-15, -1.66533454e-16, ...,
                  0.00000000e+00,  5.68434189e-14, -4.26325641e-14])
```

```
In [109]: linalg.norm(x2.real - x)
```

```
Out[109]: 3.7388984958239193e-13
```

1.4 Números aleatorios

Hay un submódulo de `numpy` para producir números (pseudo-)aleatorios.

```
In [1]: from numpy import random
```

```
In [2]: random.rand
```

```
Out[2]: <function rand>
```

```
In [26]: # numero aleatorio uniforme entre 0 y 1:
         random.rand()
```

```
Out[26]: 0.22360738532725755
```

```
In [27]: random.rand(3)
```

```
Out[27]: array([ 0.72638377,  0.22926055,  0.95029934])
```

```
In [29]: random.rand(3, 3)
```

```
Out[29]: array([[ 0.48644947,  0.76612385,  0.63281789],
                [ 0.46679283,  0.15790784,  0.30661562],
                [ 0.56408886,  0.90344024,  0.59620715]])
```

```
In [31]: random.randn(100) # distribucion normal estandar
```

```
Out[31]: array([ 1.96923414, -0.2141449 ,  0.77744524,  0.81325598,  0.92252027,
                0.1928308 ,  0.42165311,  0.45520047, -0.58336955, -0.85428152,
                0.80092534, -1.2563949 ,  1.01468299,  1.11516912, -0.02865218,
                0.18785725, -1.07156541,  1.18276602,  1.37330898,  0.48599099,
                -2.2630245 , -1.27926794,  0.07919993, -1.02950659, -0.6633516 ,
                -1.03719324,  0.16202782, -1.49051598,  0.6922923 ,  0.8307979 ,
                -0.74656157, -0.11054679,  0.30952972, -2.07747786,  0.57449484,
                0.06757248,  0.59923819, -0.23171694,  1.86531943,  0.57038927,
                1.02167588, -0.33153726,  0.26783818, -1.47657545,  1.83127475,
                -1.90195421,  0.88035848, -0.40454902, -2.10864041,  0.57904506,
                -1.13749481,  0.87165092, -0.4041457 , -0.53853878,  0.04560867,
                -0.7052234 ,  0.58162055,  0.59935702,  0.10278931, -0.26116051,
                1.36011493, -1.00785712, -0.75239444,  2.03494972,  1.61191708,
                0.80986506, -1.92907844,  0.30636403,  0.6661681 , -1.1178212 ,
                -0.25599829, -1.02935863, -0.53004217,  1.92130323, -0.24335478,
                0.19597113,  0.66874136,  0.66831047, -0.50718818,  0.39543966,
                0.33500129,  1.6545728 ,  0.68671895, -0.49427429, -2.24993033,
                -0.42802322,  0.79950429, -0.72465434,  0.12479739, -0.71969141,
                1.21309528, -1.21261034, -2.25926955, -0.96484993, -0.62112468,
                -1.16547805,  0.77298085, -1.46450965,  0.44122903,  0.58649299])
```

```
In [33]: random?
```

```
In [34]: M = random.rand(1000, 1000)
```

```
In [35]: M = M + M.T # produce matriz simetrica
```

```
In [36]: from numpy import linalg
```

```
In [37]: %time linalg.eigvalsh(M)
```


CPU times: user 440 ms, sys: 67.7 ms, total: 507 ms
Wall time: 139 ms

```
Out[37]: array([ -2.59571854e+01, -2.57299787e+01, -2.54592018e+01,
-2.49899836e+01, -2.48942731e+01, -2.46988124e+01,
-2.46797772e+01, -2.44949018e+01, -2.42906139e+01,
-2.41135869e+01, -2.39871592e+01, -2.38994431e+01,
-2.38540030e+01, -2.37959517e+01, -2.37655362e+01,
-2.36804965e+01, -2.36255161e+01, -2.33977906e+01,
-2.31621569e+01, -2.30941569e+01, -2.30088420e+01,
-2.29286029e+01, -2.28504189e+01, -2.28084933e+01,
-2.27789567e+01, -2.25754548e+01, -2.25250255e+01,
-2.24436788e+01, -2.23909176e+01, -2.22839870e+01,
-2.21472659e+01, -2.21144888e+01, -2.20553463e+01,
-2.19504286e+01, -2.19323138e+01, -2.17937363e+01,
-2.17110782e+01, -2.16580700e+01, -2.16423028e+01,
-2.15461161e+01, -2.14595710e+01, -2.13990163e+01,
-2.13684184e+01, -2.12742759e+01, -2.11878344e+01,
-2.11295253e+01, -2.10653198e+01, -2.10335375e+01,
-2.09480234e+01, -2.09167383e+01, -2.08469695e+01,
-2.07307874e+01, -2.06416286e+01, -2.05741497e+01,
-2.05250138e+01, -2.04490390e+01, -2.03405377e+01,
-2.02796617e+01, -2.02460154e+01, -2.01530022e+01,
-2.01084493e+01, -2.00757188e+01, -1.99006780e+01,
-1.98496158e+01, -1.97561349e+01, -1.96964970e+01,
-1.96656308e+01, -1.95761000e+01, -1.95206148e+01,
-1.94358784e+01, -1.93951500e+01, -1.93543071e+01,
-1.92968342e+01, -1.92615048e+01, -1.92245567e+01,
-1.92167195e+01, -1.90951318e+01, -1.90196994e+01,
-1.89254097e+01, -1.88326114e+01, -1.87988905e+01,
-1.87722206e+01, -1.86947710e+01, -1.86610355e+01,
-1.85928960e+01, -1.85828244e+01, -1.85276863e+01,
-1.84067701e+01, -1.83712463e+01, -1.83303537e+01,
-1.82568289e+01, -1.81884963e+01, -1.81501701e+01,
-1.80660149e+01, -1.79789636e+01, -1.79103198e+01,
-1.78857926e+01, -1.78109120e+01, -1.77559564e+01,
-1.76888325e+01, -1.76338487e+01, -1.75793250e+01,
-1.75627352e+01, -1.74880395e+01, -1.74488502e+01,
-1.74043985e+01, -1.73412819e+01, -1.72794186e+01,
-1.72509437e+01, -1.71702516e+01, -1.71248681e+01,
-1.70643767e+01, -1.69984289e+01, -1.69873667e+01,
-1.68437489e+01, -1.68266454e+01, -1.67952980e+01,
-1.67850035e+01, -1.67276305e+01, -1.66528077e+01,
-1.65818271e+01, -1.65237633e+01, -1.64813059e+01,
-1.64776294e+01, -1.63796952e+01, -1.63510813e+01,
-1.62763064e+01, -1.62697821e+01, -1.61704188e+01,
-1.61445620e+01, -1.60800758e+01, -1.60141007e+01,
-1.59689061e+01, -1.59384472e+01, -1.59059373e+01,
-1.58500212e+01, -1.57663899e+01, -1.57309559e+01,
-1.57168118e+01, -1.56219017e+01, -1.55764615e+01,
-1.54883244e+01, -1.54562439e+01, -1.54413983e+01,
-1.54353194e+01, -1.53611900e+01, -1.53412804e+01,
-1.52204081e+01, -1.51463658e+01, -1.51101475e+01,
-1.50884139e+01, -1.50264508e+01, -1.49679469e+01,
```

-1.49014151e+01, -1.48584081e+01, -1.47842729e+01,
 -1.47582447e+01, -1.47431217e+01, -1.46750662e+01,
 -1.46470457e+01, -1.45725317e+01, -1.45274933e+01,
 -1.44884967e+01, -1.44624752e+01, -1.43877117e+01,
 -1.43365492e+01, -1.42899282e+01, -1.42306994e+01,
 -1.41688938e+01, -1.41039756e+01, -1.40703656e+01,
 -1.40336978e+01, -1.40236085e+01, -1.39064143e+01,
 -1.38711941e+01, -1.38474939e+01, -1.38086233e+01,
 -1.37317268e+01, -1.36988139e+01, -1.36526632e+01,
 -1.36020624e+01, -1.35306068e+01, -1.34786099e+01,
 -1.34707666e+01, -1.34149486e+01, -1.33540104e+01,
 -1.33181162e+01, -1.32782382e+01, -1.32381997e+01,
 -1.31888894e+01, -1.31010430e+01, -1.30810553e+01,
 -1.30304858e+01, -1.29675313e+01, -1.29308300e+01,
 -1.29102870e+01, -1.28277364e+01, -1.27954067e+01,
 -1.27410790e+01, -1.27041295e+01, -1.26661560e+01,
 -1.26109586e+01, -1.25771604e+01, -1.25546716e+01,
 -1.24854415e+01, -1.24587167e+01, -1.24292781e+01,
 -1.23404477e+01, -1.22985216e+01, -1.22762812e+01,
 -1.22290008e+01, -1.22049958e+01, -1.21431863e+01,
 -1.20738053e+01, -1.20530760e+01, -1.20216106e+01,
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2.05702969e+01,	2.06396645e+01,	2.06879401e+01,
2.07493466e+01,	2.08532166e+01,	2.09245653e+01,
2.09929384e+01,	2.10668321e+01,	2.10826681e+01,
2.11225779e+01,	2.12289418e+01,	2.13066334e+01,
2.13867166e+01,	2.15091833e+01,	2.15234533e+01,
2.15797594e+01,	2.17294087e+01,	2.18113567e+01,

```

2.18636895e+01, 2.19642491e+01, 2.20002768e+01,
2.20518529e+01, 2.21089959e+01, 2.22828218e+01,
2.23244636e+01, 2.24295550e+01, 2.24854301e+01,
2.25656223e+01, 2.26827262e+01, 2.27531006e+01,
2.28970742e+01, 2.29598259e+01, 2.30413797e+01,
2.31260885e+01, 2.32207064e+01, 2.32732181e+01,
2.33718843e+01, 2.34205880e+01, 2.36303324e+01,
2.36669432e+01, 2.37529847e+01, 2.38455753e+01,
2.39453285e+01, 2.40423190e+01, 2.41505027e+01,
2.42205346e+01, 2.44647644e+01, 2.45482493e+01,
2.46440908e+01, 2.47100467e+01, 2.48141223e+01,
2.49162433e+01, 2.51373548e+01, 2.54662104e+01,
1.00105564e+03])

```

```

In [15]: M = np.zeros( (2, 2, 2) )
M

```

```

Out[15]: array([[[ 0.,  0.],
                  [ 0.,  0.]],

               [[ 0.,  0.],
                  [ 0.,  0.]])

```

```

In [17]: M = np.zeros( (2, 2, 2, 2) )
M

```

```

Out[17]: array([[[[ 0.,  0.],
                   [ 0.,  0.]],

                 [[ 0.,  0.],
                   [ 0.,  0.]]],

               [[[ 0.,  0.],
                   [ 0.,  0.]],

                 [[ 0.,  0.],
                   [ 0.,  0.]])])

```

```

In [18]: def f(i, j):
         return i+j

```

```

In [21]: M = np.fromfunction(f, (5, 5))
M

```

```

Out[21]: array([[ 0.,  1.,  2.,  3.,  4.],
                 [ 1.,  2.,  3.,  4.,  5.],
                 [ 2.,  3.,  4.,  5.,  6.],
                 [ 3.,  4.,  5.,  6.,  7.],
                 [ 4.,  5.,  6.,  7.,  8.]])

```

```

In [23]: def g(i, j, k):
         return i+j+k

```

```

In [25]: M = np.fromfunction(g, (5, 5, 5))
M

```

```

Out[25]: array([[ 0.,  1.,  2.,  3.,  4.],
               [ 1.,  2.,  3.,  4.,  5.],
               [ 2.,  3.,  4.,  5.,  6.],
               [ 3.,  4.,  5.,  6.,  7.],
               [ 4.,  5.,  6.,  7.,  8.]],

              [[ 1.,  2.,  3.,  4.,  5.],
               [ 2.,  3.,  4.,  5.,  6.],
               [ 3.,  4.,  5.,  6.,  7.],
               [ 4.,  5.,  6.,  7.,  8.],
               [ 5.,  6.,  7.,  8.,  9.]],

              [[ 2.,  3.,  4.,  5.,  6.],
               [ 3.,  4.,  5.,  6.,  7.],
               [ 4.,  5.,  6.,  7.,  8.],
               [ 5.,  6.,  7.,  8.,  9.],
               [ 6.,  7.,  8.,  9., 10.]],

              [[ 3.,  4.,  5.,  6.,  7.],
               [ 4.,  5.,  6.,  7.,  8.],
               [ 5.,  6.,  7.,  8.,  9.],
               [ 6.,  7.,  8.,  9., 10.],
               [ 7.,  8.,  9., 10., 11.]],

              [[ 4.,  5.,  6.,  7.,  8.],
               [ 5.,  6.,  7.,  8.,  9.],
               [ 6.,  7.,  8.,  9., 10.],
               [ 7.,  8.,  9., 10., 11.],
               [ 8.,  9., 10., 11., 12.]])

```

2 Escribir y leer datos

```
In [44]: M = np.array([1., 2, 3, 40]).reshape(2,2)
```

```
In [45]: M
```

```
Out[45]: array([[ 1.,  2.],
               [ 3., 40.]])
```

```
In [50]: np.savetxt("M.dat", M, fmt="%10.5f", header="Una matriz interesante")
```

```
In [51]: %less M.dat
```

```
In [53]: N = np.loadtxt("M.dat")
```

```
In [54]: N
```

```
Out[54]: array([[ 1.,  2.],
               [ 3., 40.]])
```

```
In [57]: np.loadtxt?
```

```
In [59]: M = np.fromfunction(f, (3, 6) )
```

```
In [60]: M
```



```

Out[60]: array([[ 0.,  1.,  2.,  3.,  4.,  5.],
               [ 1.,  2.,  3.,  4.,  5.,  6.],
               [ 2.,  3.,  4.,  5.,  6.,  7.]])

In [61]: np.savetxt("M.dat", M, fmt="%10.5f", header="Una matriz interesante")

In [64]: N = np.loadtxt("M.dat", usecols=(2, 4, 5))
        N

Out[64]: array([[ 2.,  4.,  5.],
               [ 3.,  5.,  6.],
               [ 4.,  6.,  7.]])

In [65]: x, y, z = N.T

In [66]: x

Out[66]: array([ 2.,  3.,  4.])

In [69]: x, y, z = np.loadtxt("M.dat", usecols=(2, 4, 5), unpack=True)

In [70]: y

Out[70]: array([ 4.,  5.,  6.])

```

2.1 Sustitución en cadenas

```

In [71]: dx = 0.1
        nombre_archivo = "ecn_calor_dx.dat"
        nombre_archivo

Out[71]: 'ecn_calor_dx.dat'

In [76]: dx = 0.1
        nombre_archivo = "ecn_calor_%s.dat"
        nombre_archivo

Out[76]: 'ecn_calor_%s.dat'

In [77]: nombre_archivo % dx

Out[77]: 'ecn_calor_0.1.dat'

In [83]: nombre_archivo = "ecn_calor_{0}_dx{0}.dat".format(dx)
        nombre_archivo

Out[83]: 'ecn_calor_0.1_dx0.1.dat'

In [85]: !./ecn_calor $dx

/bin/sh: ./ecn_calor: No such file or directory

In [86]: %%bash
        ls
        echo "Hola"

```

```

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raices.pyc
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test_dif_auto.pyc
test_intervalo.py
test_intervalo.pyc
Hola

In [87]: a = _

In [88]: a

Out[88]: './ecn_calor 0.1'

In [89]: a = 1

In [90]: s = "a = a y b = b"

In [94]: s1 = "Hola"
         s2 = "David"

```

```

s1 + ", " + s2
Out[94]: 'Hola, David'

In [99]: s1 + ", " + s2 + ", el valor de a es " + str(a)
Out[99]: 'Hola, David, el valor de a es 1'

In [100]: s = "3"
In [103]: int(s)
Out[103]: 3

In [104]: float(s)
Out[104]: 3.0

In [107]: dx = 0.1
          dy = 0.1
          s = "ecn_calor_dx" + str(dx) + "_dy" + str(dy)
          s

Out[107]: 'ecn_calor_dx0.1_dy0.1'

In [118]: s = "ecn_calor_%s"
In [119]: s
Out[119]: 'ecn_calor_%s'

Un nuevo operador %:

In [120]: s % dx
Out[120]: 'ecn_calor_0.1'

In [121]: dx = 0.1
          dy = 0.2

In [125]: s = "ecn_calor_dx%s_dy%s" % (dx, dy)
          s

Out[125]: 'ecn_calor_dx0.1_dy0.2'

Otra sintaxis (más nueva):

In [127]: s = "ecn_calor_dx{}_dy{}".format(dx, dy)
          s

Out[127]: 'ecn_calor_dx0.1_dy0.2'

In []:

```