NMEP HW #1

1. a. (17+ (3) = (47)

b. (3) (3) (4) = (8)

 $C = \begin{pmatrix} 0 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \end{pmatrix} = \begin{pmatrix} 2 & 1 & 1 \\ 1 & 1 & 1 \end{pmatrix}$

a. Wilspane: Ax=0

123 0

R3=R2-P3 (230)

 $R_1 = R_1 - \frac{1}{3}R_2$ $\begin{pmatrix} 1 & 0 & \frac{1}{3} & 0 \\ 0 & 6 & 5 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix}$

 $\overrightarrow{x} = \begin{pmatrix} -\frac{1}{6} \\ -\frac{1}{6} \end{pmatrix}$

Columnspace: Only 2 of thereits are knearly independent. so whom space is Span & (-1) (472)

b. (No), a linear transformation can be represented by a matrix. But fir 2 matrices AB is not necessarily equal to BA. For example, if A = [3 47 and B= (56),

2. b (lamt). AB=
$$\begin{bmatrix} \frac{12}{2} & \frac{1}{2} & \frac{66}{4} & \frac{1}{2} & \frac{12}{4} & \frac{20}{3} \end{bmatrix}$$

BA = $\begin{bmatrix} \frac{56}{12} & \frac{12}{12} & \frac{12}{2} & \frac{20}{3} & \frac{31}{3} \end{bmatrix}$

Prefore, $T_{1}(T_{2}(\nabla)) = T_{2}(T_{1}(\nabla))$ is and class sorrest.

C. (No), take this words example, If $T_{1} = \begin{bmatrix} \frac{1}{12} & \frac{1}{$

3. a (wh). =
$$\begin{pmatrix} 3 & 0 & 1 & 1 & 1 & 1 \\ 0 & 24 & 1 & 1 & 1 & 1 & 1 \end{pmatrix}$$

= $\frac{1}{12} \begin{pmatrix} 24 & 0 & 3 & 1 & 1 & 1 \\ 0 & 24 & 1 & 1 & 2 & 1 & 1 & 1 \end{pmatrix}$

= $\frac{1}{12} \begin{pmatrix} 24 & 0 & 3 & 1 & 1 & 1 \\ 0 & 24 & 1 & 1 & 2 & 1 & 1 & 1 \end{pmatrix}$

b. $A\overrightarrow{x} = \begin{pmatrix} 1 & 2 & 1 & 2 & 1 & 1 & 1 & 1 \\ 0 & 24 & 1 & 1 & 1 & 1 & 1 & 1 & 1 \end{pmatrix}$

c. $||\overrightarrow{b} - A\overrightarrow{x}||^2 = ||(\overrightarrow{a} - 3) - (-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || = ||(-4) || =$

We can have the hyperparameter & by trying out a bank of different values and then checking him well the worresponding wi fits our days (budsearch).

Intro to Numpy

What is Numpy and why do we use it?

It's an awesome python package that adds support for large, multi-dimensional arrays. Really good for vector operations, matrix operations because its super parallelized so its super fast!

Why not Python arrays?

Python arrays has certain limitations: they don't support "vectorized" operations like elementwise addition and multiplication, and the fact that they can contain objects of differing types mean that Python must store type information for every element, and must execute type dispatching code when operating on each element. This also means that very few list operations can be carried out by efficient C loops – each iteration would require type checks and other Python API bookkeeping.

Importing numpy

Functions for numerical computiing are provided by a separate module called numpy which we must import.

By convention, we import numpy using the alias np.

Once we have done this we can prefix the functions in the numpy library using the prefix np.

```
In [1]: # This is the de facto way to import NumPy. You probably don't want to
    write numpy.whatever every time
    import numpy as np
```

Numpy Arrays

NumPy arrays are the workhorse of the library. A NumPy array is essentially a bunch of data coupled with some metadata:

type: the type of objects in the array. This will typically be floating-point numbers for our purposes, but other types can be stored. The type of an array can be accessed via the dtype attribute.

shape: the dimensions of the array. This is given as a tuple, where element i of the tuple tells you how the "length" of the array in the ith dimension. For example, a 10-dimensional vector would have shape (10,), a 32-by-100 matrix would have shape (32,100), etc. The shape of an array can be accessed via the shape attribute.

Let's see some examples! There are number of ways to construct arrays. One is to pass in a Python sequence (such as list or tuple) to the np.array function:

```
In [2]: np.array([1, 2.3, -6])
Out[2]: array([ 1. , 2.3, -6. ])
```

We can also easily create ordered numerical lists!

```
In [3]: # Remember we zero index so you will actually get 0 to 6!
print(np.arange(7))
# Remember the list wont include 9
print(np.arange(3, 9))

[0 1 2 3 4 5 6]
[3 4 5 6 7 8]
```

We can also customize these list with a third paramter that specifices step size

```
In [4]: np.arange(0.0, 100.0, 10.0)
Out[4]: array([ 0., 10., 20., 30., 40., 50., 60., 70., 80., 90.])
```

To create a multi-dimensional array, we'll need to nest the sequences:

Neat!

There are also many convenience functions for constructing special arrays. Here are some that might be useful:

```
In [6]: # The identity matrix of given size
        np.eye(7)
Out[6]: array([[1., 0., 0., 0., 0., 0., 0.],
               [0., 1., 0., 0., 0., 0., 0.]
               [0., 0., 1., 0., 0., 0., 0.]
               [0., 0., 0., 1., 0., 0., 0.]
               [0., 0., 0., 0., 1., 0., 0.]
               [0., 0., 0., 0., 0., 1., 0.],
               [0., 0., 0., 0., 0., 0., 1.]])
In [7]: # A matrix with the given vector on the diagonal
        np.diag([1.1,2.2,3.3])
Out[7]: array([[1.1, 0. , 0.],
               [0., 2.2, 0.],
               [0., 0., 3.3]])
In [8]: #An array of all zeros or ones with the given shape
        np.zeros((8,4)), np.ones(3)
Out[8]: (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.]]), array([1., 1., 1.]))
In [9]: # An array with a given shape full of a specified value
        np.full((3,4), 2.1)
Out[9]: array([[2.1, 2.1, 2.1, 2.1],
               [2.1, 2.1, 2.1, 2.1],
               [2.1, 2.1, 2.1, 2.1]
```

```
In [10]:
         # A random (standard normal distribution) array with the given shape
         np.random.randn(5,6)
Out[10]: array([[-0.21012433, -0.395064 , 0.58699609, 0.21996799, -0.63239
         225,
                  0.26370822],
                [-0.46016365, -1.18521576, -0.40156712, -0.01302791,
         806,
                  0.75917215],
                [-0.1697336, 1.30926073, 0.68306209, 0.05048018, -0.25175]
         731,
                  1.06677987],
                [-0.42961748, -0.34984805, -0.7626864, -0.65096765,
                                                                      0.66512
         512,
                 -0.2647765 ],
                              0.4199001 , 0.28404666, -0.66111752, 1.26021
                [ 0.10989286,
         329,
                  0.25438349]])
```

Okay your turn! In the cell belows try and create:

A diagonal matrix with values from 1-20 (try and create this and only type two numbers!)

```
#Your answer here
In [11]:
           np.diag(np.arange(1,21))
                                                0,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
Out[11]: array([[ 1,
                            0,
                                 0,
                                      0,
                                           0,
                                                                         0,
                                                                                   0,
                                                                                        0,
                                                                                             0,
           0,
                       0,
                            0,
                                 0,
                                      0],
                                           0,
                                                                                   0,
                            2,
                                 0,
                                      0,
                                                0,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
                                                                         0,
                                                                              0,
                    [ 0,
           0,
                                      0],
                            0,
                                 0,
                                                0,
                                                                         0,
                            0,
                                           0,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
                    [ 0,
                                 3,
                                      0,
                                                                                   0,
           0,
                       0,
                            0,
                                 0,
                                      0],
                                      4,
                                           0,
                                                0,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
                                                                         0,
                                                                              0,
                                                                                   0,
                    [ 0,
                            0,
                                 0,
           0,
                                 0,
                       0,
                            0,
                                      0],
                                      0,
                                           5,
                                                0,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
                                                                         0,
                                                                                   0,
                    [ 0,
           0,
                                      0],
                       0,
                            0,
                                 0,
                                                                              0,
                    [ 0,
                            0,
                                 0,
                                      0,
                                           0,
                                                6,
                                                     0,
                                                          0,
                                                               0,
                                                                    0,
                                                                         0,
                                                                                   0,
                                                                                        0,
           0,
                       0,
                                 0,
                            0,
                                      0],
                                                0,
                                      0, 0,
                                                    7,
                                                          0,
                                                               0,
                                                                    0,
                                                                         0,
                                                                              0,
                    [ 0,
                            0,
                                 0,
                                                                                   0,
           0,
                                 0,
                                      0],
                       0,
                            0,
                                                0,
                                                                    0,
                                                               0,
                                           0,
                                                     0,
                                                          8,
                                                                         0,
                                                                              0,
                    [ 0,
                                      0,
                                                                                   0,
                                                                                        0,
```

```
Ο,
            0,
                 0,
                      0,
                           0],
                                                      9,
                                 0,
                                      0,
                                           0,
                                                 0,
                                                          0,
                                                                 0,
                                                                      0,
                 0,
                      0,
                            0,
                                                                            0,
         [ 0,
0,
            0,
                 0,
                            0],
                                                                 0,
                                 0,
                                                 0,
                                                      0, 10,
                                                                      0,
                                                                            0,
                 0,
                      0,
                            0,
                                      0,
                                            0,
         [ 0,
0,
            0,
                 0,
                            0],
                                      0,
         [ 0,
                      0,
                           0,
                                 0,
                                            0,
                                                 0,
                                                      0,
                                                           0, 11,
                                                                      0,
                                                                            0,
                 0,
0,
            0,
                                      0,
                      0,
                           0,
                                 0,
                                           0,
                                                 0,
                                                      0,
                                                           0,
                                                                 0, 12,
                                                                            0,
                                                                                      0,
         [ 0,
                 0,
                                                                                 0,
0,
            0,
                 0,
                            0],
                                      0,
                                                           0,
                                 0,
                                                      0,
                                                                 0,
                            0,
                                            0,
                                                 0,
                                                                      0, 13,
         [ 0,
                 0,
                      0,
0,
            0,
                 0,
                      0,
                            0],
                                 0,
                                      0,
                                            0,
                                                 0,
                                                      0,
                                                            0,
         [ 0,
                 0,
                      0,
                           0,
                                                                 0,
                                                                      0,
                                                                            0, 14,
0,
            0,
                 0,
                      0,
                            0],
                                                      0,
                                                           0,
                                           0,
                                                                      0,
                                 0,
                                      0,
                                                                                 0, 15,
         [ 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,
16,
            0,
                 0,
                           0],
                                 0,
                                      0,
                                           0,
                                                 0,
                           0,
                                                      0,
                                                            0,
                                                                 0,
                                                                      0,
         [ 0,
                 0,
                      0,
                                                                            0,
0,
          17,
                 0,
                      0,
                            0],
                                      0,
                                                                           0,
                                 0,
         [ 0,
                 0,
                      0,
                           0,
                                            0,
                                                 0,
                                                      0,
                                                            0,
                                                                 0,
                                                                      0,
0,
                      0,
            0, 18,
                           0,
                                 0,
                                      0,
                                           0,
                                                 0,
                                                           0,
                                                                            0,
                 0,
                      0,
                                                      0,
                                                                 0,
         [ 0,
0,
            0,
                 0, 19,
                            0],
                           0,
                                 0,
                                      0,
                                            0,
                                                 0,
                                                      0,
                                                            0,
                                                                 0,
                                                                      0,
                                                                            0,
         [ 0,
                 0,
                      0,
                                                                                 0,
0,
            0,
                 0,
                      0, 20]])
```

Okay now let's suppose we have some data in an array so we can start doing stuff with it.

```
In [12]:
         A = np.random.randn(10,5); x = np.random.randn(5)
Out[12]: array([[-2.92468280e-01, 8.79546762e-01, -1.31717614e+00,
                  1.08877423e+00, -1.13145551e+00],
                [ 1.32223145e+00, 6.74591595e-01, -6.81043247e-01,
                 -7.36051389e-01, -8.34278291e-02],
                [-4.68542890e-01, 4.42117298e-01, 1.21137837e+00,
                 -1.33352354e+00, -1.38875897e+001,
                [ 2.02622100e-01, -5.24428790e-01, -1.04213330e+00,
                  3.22529310e-01, -9.88326861e-01],
                [-1.20211711e-03, 1.02885550e+00, 5.90903166e-01,
                 -6.08259125e-01, -1.58422186e+00],
                [-3.67900439e-01, 1.46486428e+00, 9.47578215e-01,
                 -8.01376683e-01, 1.64676644e-01],
                [-4.11347485e-01, -8.57052284e-01, -1.46004739e+00,
                 -3.91742223e-01, 6.33541639e-02],
                [-2.27954001e-01, -6.15088897e-01,
                                                    8.01566650e-01,
                  7.83904053e-01, 4.73235355e-01],
                [-7.31756484e-01, 1.52676976e-01,
                                                   1.08942979e+00,
                 -4.37079083e-01, -5.59649993e-01],
                [-1.63605277e+00, -3.19578330e-01, -3.16311099e-01,
                 -2.49333922e+00, 7.77339143e-02]])
```

One useful thing that NumPy lets us do efficiently is apply the same function to every element in an array. You'll often need to e.g. exponentiate a bunch of values, but if you use a list comprehension or map with the builtin Python math functions it may be really slow. Instead just write

We can take the sum/mean/standard deviation/etc. of all the elements in an array:

```
In [14]: np.sum(x), np.mean(x), np.std(x)
Out[14]: (-0.762267691287962, -0.1524535382575924, 0.4153529425783442)
```

You can also specify an axis over which to compute the sum if you want a vector of row/column sums (again, sum here can be replaced with mean or other operations):

Linear Algebra

By now we have a pretty good idea of how data is stored and accessed within NumPy arrays. But we typically want to do something more "interesting", which for our ML purposes usually means linear algebra operations. Fortunately, numpy has good support for such routines. Let's see some examples!

```
# Transpose a matrix
In [17]:
Out[17]: array([[-2.92468280e-01, 1.32223145e+00, -4.68542890e-01,
                  2.02622100e-01, -1.20211711e-03, -3.67900439e-01,
                 -4.11347485e-01, -2.27954001e-01, -7.31756484e-01,
                 -1.63605277e+001,
                [ 8.79546762e-01, 6.74591595e-01, 4.42117298e-01,
                 -5.24428790e-01, 1.02885550e+00, 1.46486428e+00,
                 -8.57052284e-01, -6.15088897e-01,
                                                    1.52676976e-01,
                 -3.19578330e-01],
                [-1.31717614e+00, -6.81043247e-01, 1.21137837e+00,
                 -1.04213330e+00, 5.90903166e-01, 9.47578215e-01,
                 -1.46004739e+00, 8.01566650e-01, 1.08942979e+00,
                 -3.16311099e-01],
                [1.08877423e+00, -7.36051389e-01, -1.33352354e+00,
                  3.22529310e-01, -6.08259125e-01, -8.01376683e-01,
                 -3.91742223e-01, 7.83904053e-01, -4.37079083e-01,
                 -2.49333922e+00],
                [-1.13145551e+00, -8.34278291e-02, -1.38875897e+00,
                 -9.88326861e-01, -1.58422186e+00, 1.64676644e-01,
                  6.33541639e-02, 4.73235355e-01, -5.59649993e-01,
                  7.77339143e-02]])
```

Now that you're familiar with numpy feel free to check out the documentation and see what else you can do! Documentation can be found here: https://docs.scipy.org/doc/ (https://docs.scipy.org/doc/)

Exercises

Lets try out all the new numpy stuff we just learned! Even if you have experience in numpy we suggest trying these out.

1) Create a vector of size 10 containing zeros

```
In [18]: ## FILL IN YOUR ANSWER HERE ##
v = np.zeros(10)
v

Out[18]: array([0., 0., 0., 0., 0., 0., 0., 0., 0.])
```

2) Now change the fith value to be 5

```
In [20]: ## FILL IN YOUR ANSWER HERE ##
v[4] = 5
v
Out[20]: array([0., 0., 0., 0., 5., 0., 0., 0., 0., 0.])
```

3) Create a vector with values ranging from 10 to 49

4) Reverse the previous vector (first element becomes last)

5) Create a 3x3 matrix with values ranging from 0 to 8. Create a 1D array first and then resshape it. </br>

6) Create a 3x3x3 array with random values

7) Create a random array and find the sum, mean, and standard deviation

```
In [26]: ## FILL IN YOUR ANSWER HERE ##
v = np.random.randn(100)
np.sum(v), np.mean(v), np.std(v)

Out[26]: (0.20254356081693747, 0.0020254356081693746, 1.0827826632876085)
In []:
```

Linear Regression - Gaussian Orbits

Imports needed in this notebook: numpy (as np), matplotlib.pyplot (as plt), from sklearn: LinearRegression, ElasticNet, and mean_squared_error. Search up documentation if you have issues with any of the scikit learn things. This first part will be a warmup with numpy, linear regression, and scikit learn.

```
In [3]: ### YOUR IMPORT CODE HERE
   import numpy as np
   import matplotlib.pyplot as plt
   from sklearn.linear_model import LinearRegression, ElasticNet
   from sklearn.metrics import mean_squared_error
```

First, we'll generate some random data and find the line of best fit.

```
In [7]: num_data = 20

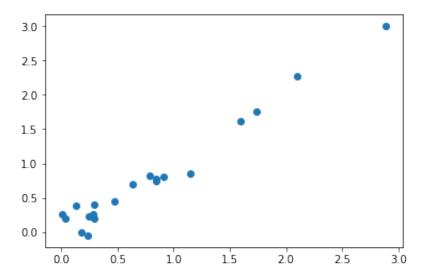
def gen_data(n):
    x = []
    y = []

    for i in range(n):
        rand = abs(np.random.randn())
        x += [rand]
        y += [.15 * np.random.randn() + rand]

    plt.scatter(x, y)
    plt.show()

    return (x, y)

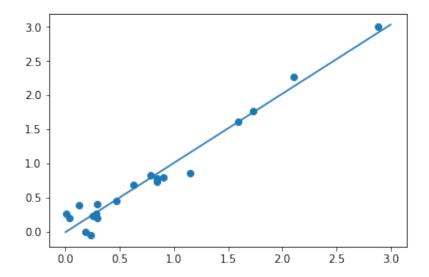
data = gen_data(num_data)
    data = [(data[0][i], data[1][i]) for i in range(num_data)]
    data
```



```
Out[7]:
        [(0.8463782847314022, 0.7362085678615494),
         (0.7870448153101179, 0.8252617577179918),
         (0.12874811072226117, 0.388945908839599),
         (0.8441908955035079, 0.7811857947591122),
         (0.277997439039963, 0.26757260272311256),
         (1.149828611799792, 0.8579499967276797),
         (0.9062900503387852, 0.8001301680305749),
         (0.46880128078087707, 0.45253152660085294),
         (1.5940586932725254, 1.6134183106313167),
         (0.009976240390721394, 0.2670116990253635),
         (1.7355415858677445, 1.7624280826362329),
         (0.6315276383656515, 0.6909427154520842),
         (0.294433815590066, 0.20399015461893544),
         (2.886809868918759, 3.0012619737982424),
         (0.18209952293050177, -0.002718136538349153),
         (2.1008826641350455, 2.2749239682811884),
         (0.2306681431508596, -0.05234509877551913),
         (0.24794407292381804, 0.2333180954596053),
         (0.29437782244312527, 0.4037965231652011),
         (0.034698550152521435, 0.19882754911689532)
```

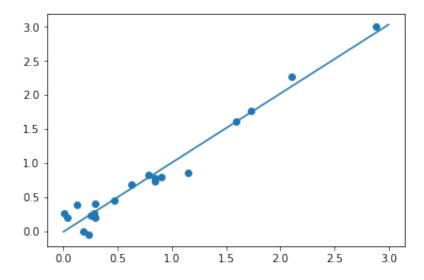
Numpy Linear Regression

Now, data is a list of tuples of data. Perform linear regression to find the line of best fit, using only numpy. Minimize squared loss. Use matplotlib to plot your line of best fit.



Sklearn Linear Regression

Now, do the same thing, except using Scikit Learn. I recommend reading some documentation, specifically here: https://scikit-learn.org/stable/modules/generated/sklearn.linear_model.LinearRegression.html)



Gaussian Orbits

In this homework, we will use linear regression methods in order to determine the orbits of heavenly bodies.

Background

In 1801 the minor planet Ceres was first observed for a period of 40 days before moving behind the sun. To predict the location of Ceres astronomers would have to solve complicated non-linear differential equations, quite a task in an era before computers or calculators. However, Carl Friedrich Gauss had another idea. By single handedly developing the theory of least squares and linear regression and applying it to the problem of finding Ceres, Gauss managed to accurately predict the location of the minor planet nearly a year after it's last sighting.

In this problem we likewise attempt to predict the orbit of a "planet" and in the process "derive" the formula for an ellipse, the shape of orbits of heavenly bodies.

1. Generate Data

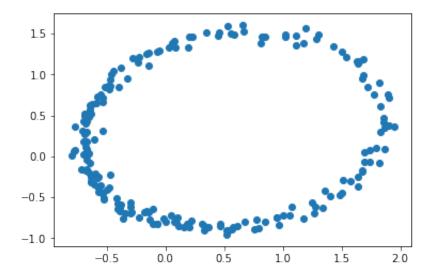
The idea here is we generate data in the shape of an ellipse. To do this we use the formula of an ellipse in polar coordinates:

$$r = \frac{ep}{1 - e\cos\theta}$$

where e is the eccentricity and p is the distance from the minor axis to the directrix (read "length"). In addition, we add random noise to the data.

We will then try to fit curves to our synthetic dataset.

```
In [57]: def gen data(e, p, o):
             theta = np.linspace(0,2*np.pi, 200)
             # Ellipse with eccentricity e
             # Axis "length" p
             # Offset by .5 angularly
             r = e*p/(1-e*np.cos(theta - o))
             # transform to cartesian
             x = r * np.cos(theta)
             y = r * np.sin(theta)
             # Add noise
             x += np.random.randn(x.shape[0]) / 20
             y += np.random.randn(y.shape[0]) / 20
             # plot
             plt.scatter(x, y)
             plt.show()
             # saving
             np.save('x.npy', x)
             np.save('y.npy', y)
             return x, y
         x, y = gen_data(.5, 2, .5)
```



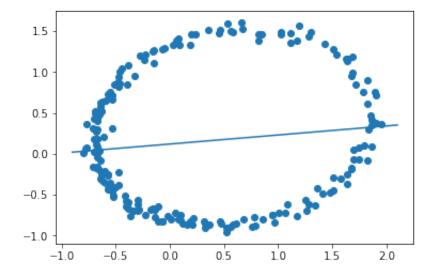
2. Use sklearn's LinearRegression

Try to fit a LinearRegression model to x and y (let x be the independent variable and y be the dependent variable). Print out the mean_squared_error you get and plot both x, y (scatter plot), and the predicted orbit (line plot).

```
In [59]: ### YOUR CODE HERE
model = LinearRegression()
model.fit(x.reshape(-1, 1), y)

plt.scatter(x, y)
    line_x_data = np.linspace(-0.9, 2.1, 1000)
    plt.plot(line_x_data, model.coef_[0] * line_x_data + model.intercept_)
    plt.show()

print("Mean squared error: " + str(mean_squared_error(y, model.coef_[0] * x + model.intercept_)))
```



Mean squared error: 0.6559930382275464

This is not the best approach for our data. Please explain why below.

The data is not linear! In other words, the "line of best fit" is not a straight line in the Cartesian plane, but rather an ellipse.

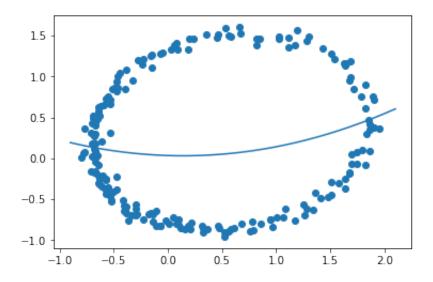
3. Experimentation time!

Try adding new features to your linear model by manipulating x! For example, try adding a quadratic term, x^2 or a root term like \sqrt{x} . Print out the MSE of your model and plot both x, y (scatter plot), and the predicted orbit (line plot). This time, your model should take in an expanded set of features and predict y.

Hint: np.vstack may be useful here.

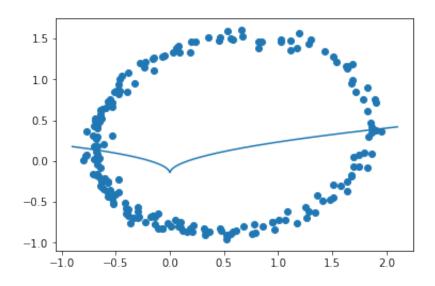
```
In [77]:
         ### YOUR CODE HERE
         print("Adding x^2 term")
         features = np.hstack((np.square(x).reshape(-1, 1), x.reshape(-1, 1)))
         model = LinearRegression()
         model.fit(features, y)
         plt.scatter(x, y)
         line_x_data = np.linspace(-0.9, 2.1, 1000)
         line_y_data = np.dot(np.hstack((np.square(line_x_data).reshape(-1, 1),
         line x data.reshape(-1, 1))), model.coef ) + model.intercept
         plt.plot(line x data, line y data)
         plt.show()
         y pred = np.dot(features, model.coef ) + model.intercept
         print("Mean squared error: " + str(mean squared error(y, y pred)))
         print("\n\nAdding sqrt(|x|) term")
         features = np.hstack((np.sqrt(np.abs(x)).reshape(-1, 1), x.reshape(-1,
         1)))
         model = LinearRegression()
         #print(features)
         model.fit(features, y)
         plt.scatter(x, y)
         line x data = np.linspace(-0.9, 2.1, 1000)
         line y data = np.dot(np.hstack((np.sqrt(np.abs(line x data)).reshape(-
         1, 1), line x data.reshape(-1, 1))), model.coef) + model.intercept
         plt.plot(line x data, line y data)
         plt.show()
         y pred = np.dot(features, model.coef ) + model.intercept
         print("Mean squared error: " + str(mean squared error(y, y pred)))
```

Adding x^2 term



Mean squared error: 0.6486438351011222

Adding sqrt(|x|) term



Mean squared error: 0.6492861409394708

This is still not the best idea, please explain:

The data clearly doesn't follow neither a quadratic curve nor a sqrt(x) curve.

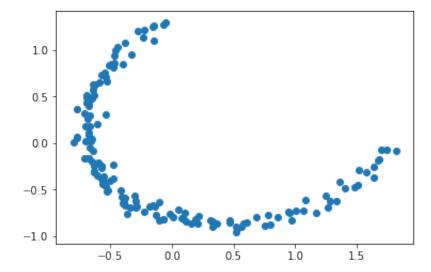
4. Plane Curves

As you've probably figured out, the above two methods are pretty crap at predicting orbits. What we really need to do is predict a curve in the plane. First, let's erase some of the data so what we're doing is actually a challenge. Just run the code in the next box:

```
In [78]: # Create a mask where x < 0 or y < 0
def mask():
    global x
    global y

    mask = (x < 0) + (y < 0)
    x = x[mask]
    y = y[mask]

# plot erased data
    plt.scatter(x, y)
    plt.show()</pre>
```



Now the most general form of a plane curve is

$$f(x, y) = 0$$

In order to simplify our lives a bit, let's restrict this to something of the form:

$$ax^2 + bxy + cy^2 + dx + ey + f = 0$$

You may recognize this as the general form of a conic! Let's take our data and try to predict the best possible coefficients here using least squares. This way, these coefficients should give the best possible approximation to the orbit. Print your predicted coefficients.

Hint 1: Think about the features you need. (6 total)

Hint 2: Use the normal equation instead of sklearn.

Hint 3: This is going to fail, why? a=b=c=d=e=f=0 works

```
In [94]: ### YOUR CODE HERE
    feature_1 = np.square(x).reshape(-1, 1)
    feature_2 = (x * y).reshape(-1, 1)
    feature_3 = np.square(y).reshape(-1, 1)
    feature_4 = x.reshape(-1, 1)
    feature_5 = y.reshape(-1, 1)
    feature_6 = np.ones(len(x)).reshape(-1, 1)
    features = np.hstack((np.hstack((np.hstack((np.hstack((feature_1, feature_2)), feature_3)), feature_4)), feature_5)), feature_6))
    b = np.zeros(len(x))
    x_hat = np.dot(np.matmul(np.linalg.inv(np.matmul(features.T, features)), features.T), b)
    print(x_hat)
```

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

5. Reformulation

The above should fail for a very trivial (pun intended) reason. The reason is that if we simply set all the coefficients to zero, we get a perfect solution! We can see this in the normal equations:

$$(A^T A)^{-1} A^T b = x$$

but $b = \vec{0}$ in our case, thus $x = \vec{0}$ trivially.

How do we get around this? One thing we can do is to not have $b = \vec{0}$. To do this, let us modify the general form of a plane curve a bit:

$$f(x, y) + 1 = 1$$

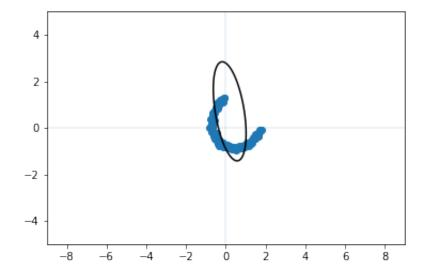
Now our restricted plane curve will be of the form

$$ax^2 + bxy + cy^2 + dx + ey + f + 1 = 1$$

Is this just an aesthetic change? or will this actually help? Code it up and find out! Plot your model using the handy dandy plot conic function

```
In [95]: # This function should help you plot your ellipses:
         def plot_conic(coeff):
             params
             _____
             coeff : array[6] floats
                 Array of 6 floats, corresponding to
                 a, b, c, d, e, and f respectively
                 in the equation above
             xv = np.linspace(-9, 9, 400)
             yv = np.linspace(-5, 5, 400)
             xv, yv = np.meshgrid(xv, yv)
             def axes():
                 plt.axhline(0, alpha=.1)
                 plt.axvline(0, alpha=.1)
             axes()
             plt.contour(xv, yv, xv*xv*coeff[0] + xv*yv*coeff[1] + yv*yv*coeff[
         2] + xv*coeff[3] + yv*coeff[4] + coeff[5], [0], colors='k')
             plt.scatter(x,y)
             plt.show()
```

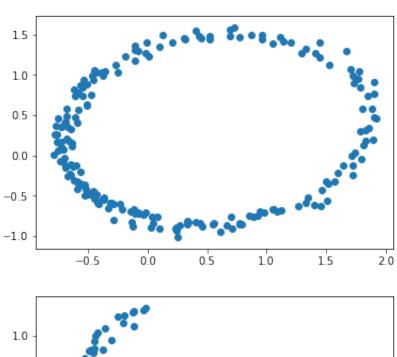
```
In [100]: ### YOUR CODE HERE
b = np.ones(len(x))
x_hat = np.dot(np.matmul(np.linalg.inv(np.matmul(features.T, features)
), features.T), b)
x_hat[5] -= 1
plot_conic(x_hat)
```

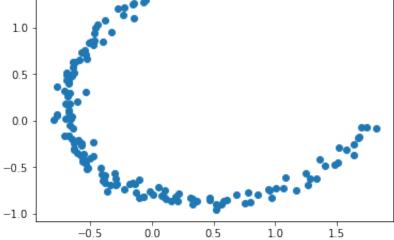


6. Ridge

So, reformulating the problem might have worked, but more than likely it didn't work too well. Here's some code to generate new data. Try running the above model multiple times on different data. More than likely most of them will look terrible.

```
In [102]: # Regenerate data
gen_data(.5, 2, .5)
mask()
```





The problem here is that our method is too unstable. It turns out the Ridge Regression as a regularizer can reduce numerical instability and constrain under-constrained problems. (The math homework with the Ridge Derivation walks you through why this is the case)

The closed form solution for Ridge Regression is the following:

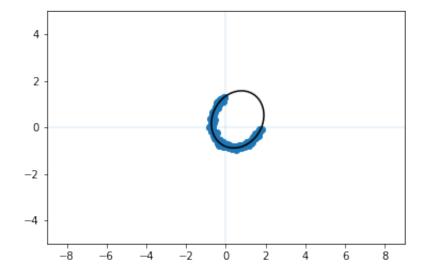
$$w_{\text{RIDGE}}^* = (X^T X + \lambda I)^{-1} X^T y$$

Rewrite the regression from above using ridge regression (try using $\lambda=1$) and see how well it does. Plot out the model using plot conic. Compare the results with the previous method.

Hint: Use the regenerate data block to try new data

Hint: There is really only one extra term between this question and the previous

```
In [103]: ### YOUR CODE HERE
b = np.ones(len(x))
ridge_lambda = 1
x_hat = np.dot(np.matmul(np.linalg.inv(np.matmul(features.T, features)
+ ridge_lambda * np.eye(features.shape[1])), features.T), b)
x_hat[5] -= 1
plot_conic(x_hat)
```



7. "Deriving" an Ellipse

LASSO regularization is a *sparse feature selector* in the sense that it zeros out "useless" features and keeps relevant features. It's a good way to reduce the number of features you have to use.

In this case we're going to pretend we don't know what form the equation of an ellipse takes. We can add random monomials to form a guess:

$$ax^{2} + bxy + cy^{2} + dx + ey + f + gx^{3} + hy^{3} + jx^{2}y + \dots + 1 = 1$$

The idea here is that if we use LASSO regression on the above equation, the terms irrelevant to an ellipse will "zero out" and the quadratic and lower terms won't! Try this out, and print out the coefficients. No gurantees this will works 100%:), but you should find that all coefficients greater than quadratic zero out.

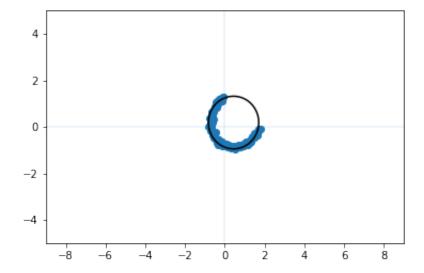
Hint: We want to keep the ridge regularization to maintain numerical stability. So we need a combined Ridge and LASSO regression. This model is called ElasticNet from sklearn. Use that model.

Hint: You might have to play around with the parameters a bit. I used these 11_ratio=.23, alpha=.01 to produce some pretty good results

```
In [127]: ### YOUR CODE HERE
          feature 7 = np.power(x, 3).reshape(-1, 1)
          feature 8 = np.power(y, 3).reshape(-1, 1)
          feature 9 = (np.power(x, 2) * y).reshape(-1, 1)
          feature 10 = (x * np.power(y, 2)).reshape(-1, 1)
          new features = np.hstack((np.hstack((np.hstack((features, f
          eature 7)), feature 8)), feature 9)), feature 10))
          net = ElasticNet(11 ratio=0.23, alpha=0.01, fit intercept=False)
          net.fit(new features, b)
          coeff = net.coef [:]
          print(coeff)
          [ 0.05565683 -0.
                                 0.06956352 - 0.04951306 - 0.02564773
                                                                      0.9241
          4928
                     -0. -0.
          -0.
                                              0.
                                                        1
```

8. Evaluate this model!

Run this code block below. This code block assumes that you have an array called <code>coeff</code> which has 10 elements.



As demonstrated above, ridge regression can help overcome numerical instability and generalization issues that ordinary least squares (OLS) can fall short to.