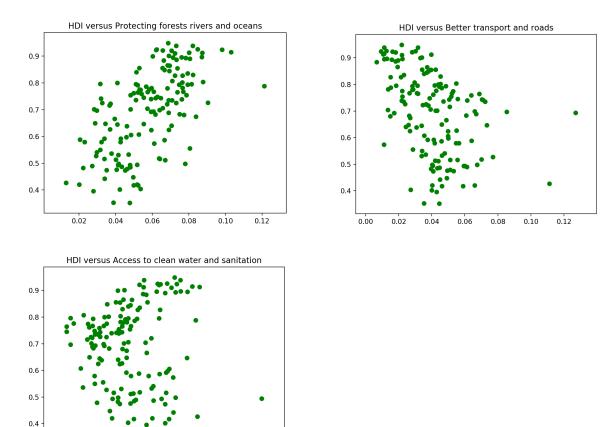
3b

Most positively correlated: Protecting forests rivers and oceans 0.613458756271

Most negatively correlated: Better transport and roads -0.439633638622

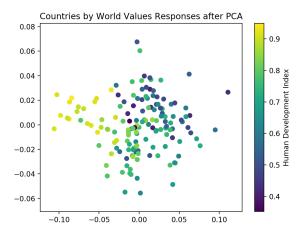
Least correlated: Access to clean water and sanitation: -0.018169084456

3c



We can observe positive correlation, negative correlation and almost no correlation in the three figures, respectively.

0.14



3e

```
Ridge Regression
RMSE: 0.12303337350607803
Coefficients
[[0.80823467 - 0.74985758 - 0.17800015 - 1.28408103 - 0.66293176 -
0.82203172
   0.73733884 - 0.92891581 - 0.82049672 0.39614952 2.0708291 -
0.06718981
   0.48310656  0.72671425  0.42921192  -0.13808023]]
3f
Lasso Regression
RMSE: 0.12602242808947528
Coefficients
[ 0.1590192 -0.72844929 -0.
                                   -0.85945074 -0.66274144 -
0.02556703
  0.33904781 -0.29897158 -0.
                                   0.
                                           3.48536375 0.
0.
  0.87057995 0.32897045 -0.
```

3g

Lasso weights have 6 zeros, while Ridge have none.

3h

Average the outputs of its k nearest nerghbors. Can also be weighted by inverse distance.

3i

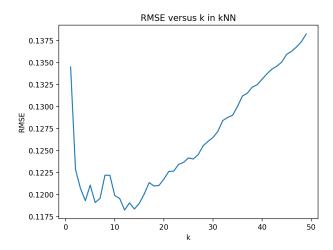
```
The closest countries are:
        United States
45
90
              Ireland
61
       United Kingdom
37
              Belgium
108
              Finland
69
                Malta
132
              Austria
110
               France
The distances are:
            0.01765082
                        0.02157067 0.03010195
0.
0.0302092
            0.03349605
   0.03583628 0.0389719
```

3j

k Nearest Neighbors Regression

RMSE: 0.11824589460776892

Best k: 12



3k

When k is large, we sampled more before making a decision, this will lead to small variance. So large k, small variance, small k, large variance.

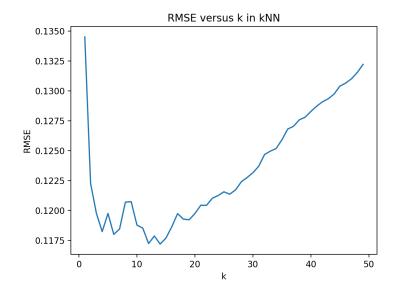
When k is large, we tend to include more samples from other classes, this will bring in bias. So large k, large bias, small k, small bias.

So we will have a sweet spot where the sum of bias and variance is the smallest.

 $\ensuremath{\mathsf{k}}$ Nearest Neighbors Regression: inverse distance weighted

RMSE: 0.1171925270311745

Best k: 14



3m

```
45
        United States
90
              Ireland
61
       United Kingdom
108
              Finland
37
              Belgium
69
                Malta
110
               France
132
              Austria
```

Compare to unscaled result, Finland and Belgium exchange positions, and France and Austria exchanged positions.

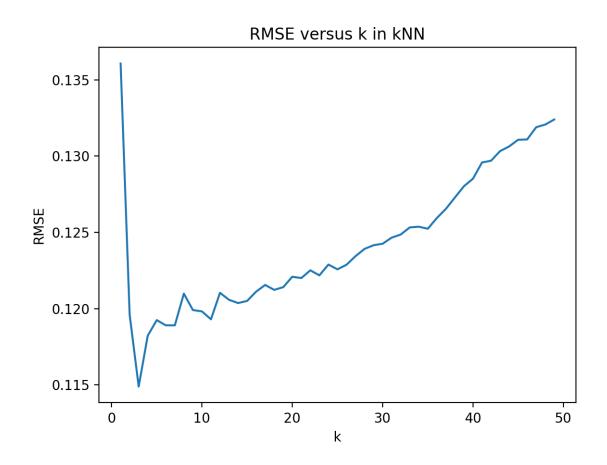
3n

k Nearest Neighbors Regression: with weights, with scaling

RMSE: 0.11488547357936414

Best k: 3

Grid search should be performed again every time we changed the setting.



3p

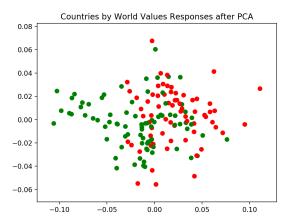
Test predictions:

- [[0.5109997]
- [0.61570753]
- [0.68331113]
- [0.67852229]
- [0.7718753]
- [0.60419024]
- [0.65159897]
- [0.67945579]
- [0.73105346]
- [0.8053592]
- [0.67938174]
- [0.66520881]
- [0.7793257]
- [0.70584762]
- [0.66568196]
- [0.56691101]
- [0.69252235]
- [0.03232233
- [0.72706267]
- [0.70498736]
- [0.77513888]
- [0.63384268]
- [0.65089538]
- [0.39709827]
- [0.73442026]
- [0.7583716]
- [0.71157266]
- [0.44080564]
- [0.90460366]
- [0.62818537]
- [0.65208228]
- [0.6630599]
- [0.86539214]
- [0.00333214]
- [0.72870152] [0.90989302]
- [0.0000000
- [0.924362] [0.8432591]
- [0.49853809]
- [0.80810577]]

3q

At least we are guaranteed a 1/k accuracy. Since the largest class should includes more than N/k samples.

3r



3s

A linear boundary cannot easily separate the two classes.

```
SVM Classification(kernel='linear', C=48.0)
Accuracy: 0.75
3u
SVM Classification(kernel='linear', PCA_k=8, C=1.0)
Accuracy: 0.77027027027 (improve a little)
3v
SVM Classification(kernel='rbf', C=73, gama=0.097)
Accuracy: 0.743243243243
3w
k Nearest Neighbors Classification(weights='distance',
k=4)
Accuracy: 0.763513513514
k Nearest Neighbors Classification(weights='distance',
k=4, scaled)
```

Accuracy: 0.77027027027(scaling helped a little)

3y

If we are given the target locations and the distance from each target for each sensor, then for each sensor, find its k nearest neighbors, and do a weighted average of the locations of the k neighbors.

The problem with this is that, if all targets are on one side, and all sensors are on the other side, then the prediction by kNN will be that all sensors are amid the targets, which is not correct.

3z

I learned that kNN can also be used for regression, and in some cases works better than ridge and lasso.

For SVM classification, 'rbf' is not always better than 'linear'.

Scaling and standardization is important.