In [1]: import pandas as pd
import numpy as np
import sklearn as sk
import scipy as scp
import xgboost as xgb
import matplotlib.pyplot as plt

In [2]: train = pd.read_csv("train.csv")
test = pd.read_csv("test.csv")

In [3]: train.head()

Out[3]:

	ACTION	RESOURCE	MGR_ID	ROLE_ROLLUP_1	ROLE_ROLLUP_2	ROLE_DEPTN/
0	1	39353	85475	117961	118300	123472
1	1	17183	1540	117961	118343	123125
2	1	36724	14457	118219	118220	117884
3	1	36135	5396	117961	118343	119993
4	1	42680	5905	117929	117930	119569

In [4]: test.head()

Out[4]:

	id	RESOURCE	MGR_ID	ROLE_ROLLUP_1	ROLE_ROLLUP_2	ROLE_DEPTNAME	F
0	1	78766	72734	118079	118080	117878	1
1	2	40644	4378	117961	118327	118507	1
2	3	75443	2395	117961	118300	119488	1
3	4	43219	19986	117961	118225	118403	1
4	5	42093	50015	117961	118343	119598	1

In [5]: train['ACTION']

Out[5]:	0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
	32739 32740 32741 32742 32743 32744 32745 32746 32747 32748 32750 32751 32752 32753 32754 32755 32756 32757 32758 32759 32760 32761 32762 32763 32763 32764	

32765

1

```
32766
         32767
                  1
         32768
                  1
         Name: ACTION, dtype: int64
 In [6]: from sklearn.linear model import LogisticRegressionCV
 In [7]: Cs = [0.0005, 0.001, 0.01, 0.1, 1, 3, 5, 10, 20, 50]
         lq = LogisticRegressionCV()
 In [8]: tr data = train.values
         y = tr data[:, 0]
         x = tr data[:, 1:]
         print tr data.shape
         print y.shape
         print x.shape
         lg.fit(x, y)
         (32769, 10)
         (32769,)
         (32769, 9)
 Out[8]: LogisticRegressionCV(Cs=10, class weight=None, cv=None, dual=False,
                    fit intercept=True, intercept scaling=1.0, max iter=100,
                    multi class='ovr', n jobs=1, penalty='l2', random state=No
         ne,
                    refit=True, scoring=None, solver='lbfgs', tol=0.0001, verb
         ose=0)
 In [9]: te_data = test.values[:, 1:]
         print te data.shape
         lgbase pred = lg.predict(te data)
         id = test.values[:, 0]
         print id[0:10]
         (58921, 9)
         [1 2 3 4 5 6 7 8 9 10]
In [10]: | lgbase pred.shape
Out[10]: (58921,)
In [11]: | id.shape
Out[11]: (58921,)
In [12]: lgbase sub = {'id':id, 'ACTION': lgbase pred}
         lgbase sub = pd.DataFrame(data = lgbase sub, columns=['id',
          'ACTION'])
In [13]: | lgbase sub.to csv('lgbase pred', index = False)
```

```
In [14]: | lgbase_sub.head()
Out[14]:
               ACTION
            1
               1
            2
               1
            3
          2
               1
          3
            4
               1
            5
               1
In [15]:
         count0 = 0
          count1 = 0
          for i in lgbase pred:
              if i == 0:count0 +=1
              if i == 1: count1 += 1
In [16]: count0
Out[16]: 0
In [17]: count1
```

As you can see, the base logistic regression model returned a prediction of all "1's". After submitting to Kaggle, we got a score of .5

Time to try one hot encoding.

Out[17]: 58921

```
In [18]: from sklearn import preprocessing
In [19]: encoder = preprocessing.OneHotEncoder()
    onehot = np.vstack((x, te_data))
    onehot = encoder.fit_transform(onehot)

In [20]: onehot_train = onehot[0:tr_data.shape[0], :]
    onehot_test = onehot[tr_data.shape[0]:, :]

In [21]: onehot_train.shape
Out[21]: (32769, 16961)
In [22]: onehot_test.shape
Out[22]: (58921, 16961)
```

Now logistic regression on this data.

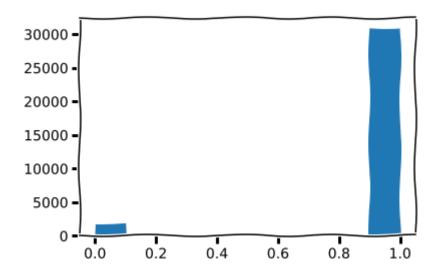
We got a 0.6375, which is a significant jump. Now lets start boosting.

```
In [28]: | from sklearn.tree import DecisionTreeRegressor
         tree = DecisionTreeRegressor()
In [29]: tree.fit(onehot train, y)
Out[29]: DecisionTreeRegressor(criterion='mse', max depth=None, max features=N
         one,
                    max leaf nodes=None, min impurity split=1e-07,
                    min samples leaf=1, min samples split=2,
                    min weight fraction leaf=0.0, presort=False, random state=
         None,
                    splitter='best')
In [30]: tree preds = tree.predict(onehot test)
         get submission(tree preds, 'tree preds')
In [31]: #how kaggle computes a score
         from sklearn import metrics
         def evaluate(preds, true labels):
             fpr, tpr, thresholds = metrics.roc curve(true labels, preds, pos
         label=1)
             auc = metrics.auc(fpr,tpr)
             return auc
```

I read up on Paul Duan's methodology. He recommends removing role 1 and role 2, due to hig variance created by these. Besides that, he doesn't worry too much about feature selection, but says feature extraction is more important. He combines different models, each using a different feature set.

```
In [32]: #first thing first, look at the data
plt.xkcd() #makes graphs look better
plt.hist(train['ACTION'])
plt.show()
```

/home/justin/anaconda2/lib/python2.7/site-packages/matplotlib/font_ma nager.py:1297: UserWarning: findfont: Font family [u'xkcd', u'Humor S ans', u'Comic Sans MS'] not found. Falling back to DejaVu Sans (prop.get family(), self.defaultFamily[fontext]))



It appears as if most predictions should be a 1. However, our first submission shows that the test set has many more zeroes than the train set.

Get rid of bad columns.

```
In [33]: del train['ROLE_ROLLUP_1']
  del train['ROLE_ROLLUP_2']
  train.head()
```

Out[33]:

	ACTION	RESOURCE	MGR_ID	ROLE_DEPTNAME	ROLE_TITLE	ROLE_FAMILY_DE
0	1	39353	85475	123472	117905	117906
1	1	17183	1540	123125	118536	118536
2	1	36724	14457	117884	117879	267952
3	1	36135	5396	119993	118321	240983
4	1	42680	5905	119569	119323	123932

```
In [34]: del test['ROLE_ROLLUP_1']
    del test['ROLE_ROLLUP_2']
    test.head()
```

Out[34]:

	id	RESOURCE	MGR_ID	ROLE_DEPTNAME	ROLE_TITLE	ROLE_FAMILY_DESC	R
0	1	78766	72734	117878	117879	118177	19
1	2	40644	4378	118507	118863	122008	11
2	3	75443	2395	119488	118172	301534	24
3	4	43219	19986	118403	120773	136187	11
4	5	42093	50015	119598	118422	300136	11

We will assume these new datasets are better bases for classification.

```
In [35]:
         #From Paul Duan
         def create tuples(X):
             cols = []
             for i in range(X.shape[1]):
                  for j in range(i, X.shape[1]):
                      cols.append(X[:, i] + X[:, j]*3571)
              return np.hstack((X, np.vstack(cols).T))
         def create_triples(X):
             cols = []
             for i in range(X.shape[1]):
                  for j in range(i, X.shape[1]):
                      for k in range(j, X.shape[1]):
                          cols.append(X[:, i]*3461 + X[:, j]*5483 + X[:, k])
             return np.hstack((X, np.vstack(cols).T))
         #mine
         def onehot(x, y):
             enc = preprocessing.OneHotEncoder()
             enc.fit(np.vstack((x, y)))
             return enc.transform(x), enc.transform(y)
In [36]: | #try to make a tuple
```

Out[38]: (32769, 7)

```
In [39]: | tr x = tuple x[0:tr x.shape[0], :]
          te x = tuple x[tr x.shape[0]:, :]
          print tr x.shape
          print te x.shape
          (32769, 35)
          (58921, 35)
In [40]:
         #so we will use the smaller dataset
          enc = sk.preprocessing.OneHotEncoder()
          y = train.values[:, 0]
          tr x = train.values[:, 1:]
          test x = test.values[:, 1:]
          enc.fit(np.vstack((tr x, test x)))
          tr x = enc.transform(tr x)
          test x fin = enc.transform(test x)
In [41]: | \text{#tuptrain } x = \text{tuple } x[0:\text{tr } x.\text{shape}[0], : ]
          #tuptest x = tuple x[tr x.shape[0]:, :]
          seed = 42
In [42]:
          test size = .20
          #y frame = train['ACTION']
          #x frame =train
          #del x frame['ACTION']
          \#x frame = enc.transform(x frame)
          print type(tr_x)
          x train, x test, y train, y test = sk.model selection.train test spli
          t(tr x, y, test size=test size, random state=seed)
          type(x_train)
         <class 'scipy.sparse.csr.csr matrix'>
Out[42]: scipy.sparse.csr.csr_matrix
In [43]:
         \#x train = np.ndarray(x train)
          #print x train
          \#x train = np.hstack((x train, (np.ones(x train.shape[0], 1))))
          #x test = scp.hstack((x test, scp.sparse.csr matrix(np.ones(x train.s
          hape[0], 1))))
          \#x \ test = np.ndarray(x \ test)
          #boosting
          xgbmodel = xgb.XGBClassifier()
          xgbmodel.fit(x_train,y_train)
Out[43]: XGBClassifier(base score=0.5, colsample bylevel=1, colsample bytree=
          1,
                 gamma=0, learning rate=0.1, max delta step=0, max depth=3,
                 min child weight=1, missing=None, n estimators=100, nthread=-
         1.
                 objective='binary:logistic', reg_alpha=0, reg_lambda=1,
                 scale pos weight=1, seed=0, silent=True, subsample=1)
```

In [44]:

print x train.shape

```
print x test.shape
         print y train.shape
         (26215, 16648)
         (6554, 16648)
         (26215,)
In [45]:
         coo = x test.tocoo(copy=False)
         x test df = pd.DataFrame({'index':coo.row, 'col':coo.col,
         'data':coo.data})[['index', 'col', 'data']].sort_values(['index', 'co
         l']).reset index(drop=True)
         #print x test1
         x train.shape
Out[45]: (26215, 16648)
In [46]: xgb preds = xgbmodel.predict(x test)
In [47]: xgb preds.shape
Out[47]: (6554,)
```

Basically, XGBoost drops a few columns because they are all zeroes. This breaks the model. This sucks, as any model is basically useless without onehot encoding. Supposedly, the developers fixed this in the newest version. However, I could not install that version. (Trust me, I tried before I gave up and used pip, which stil took significant troubleshooting.)

If this worked, I could generate a score using the returned xgb_preds and y_test values. (The score generation algorithm is available on kaggle.) Then I could manipulate the XGBoost model to try to get a better score, and submit my best scoring model. Unfortunately, the developers of XGBoost had a different vision in mind when they built their software. A vision of hopelessness and despair.

So I reinstalled and now it works. Time to start boosting for real

```
In [48]: #test accuracy
evaluate(xgb_preds, y_test)
Out[48]: 0.50804924055707812
```

This is actually worse. Time to start playing with the model.

```
In [49]: #smaller trees, less estimators
    xgb2 = xgb.XGBClassifier(max_depth = 2, learning_rate = 0.5, n_estima
    tors=50, silent=True)
    xgb2.fit(x_train, y_train)
    preds = xgb2.predict(x_test)
    evaluate(y_test, preds)
```

Out[49]: 0.87241748166259181

If this is correct, that is an insane jump.

```
In [50]: #Let's try with stumps
    xgb_stump = xgb.XGBClassifier(max_depth = 1, learning_rate= 0.5, n_es
    timators=50)
    xgb_stump.fit(x_train, y_train)
    preds = xgb_stump.predict(x_test)
    print evaluate(y_test, preds)
    print preds

0.90127785059
[1 1 1 ..., 1 1 1]
```

So this is really good. Like I don't believe it. Time to run Cross Validation.

Model 1: 0.516098481114 run time: 0: 10.6450340748 s

suck. Funny how the difference is so massive.

```
In [51]:
         #CV
         import time
         def cross val(xqbs):
             start = time.time()
             expected err = [0]*len(xgbs);
             expected err stump = 0;
             for i in range(0, 10):
                 x_train, x_test, y_train, y_test = sk.model_selection.train_t
         est split(tr x, y, test size=test size, random state=seed)
                 for i in range(0, len(xgbs)):
                     xgbs[i].fit(x train, y train)
                                                       #stump
                     expected err[i] += evaluate(xgbs[i].predict(x test), y te
         st)
             for i in range(0, len(xgbs)):
                 expected err[i] = expected err[i]/10
                 print 'Model ', i, ": ", expected_err[i]
             run time = time.time()-start
             print 'run time: ', int(run time)/60,':',run time%60, 's'
         cross val([xgb2, xgb stump])
         Model 0: 0.510678426914
```

As we can see, the previous 2 models got lucky when they presented us with high scores. In fact, they both

We also have a runtime. I will try the same code using the gpu plugin, to see if we get faster results.

```
params = {'max depth':5, 'learning rate':0.3, 'n estimators':50, 'obj
In [52]:
         ective':'binary:logistic', 'updater':'grow_gpu'}
         \#xg train = np.hstack([tr x, y])
         \#label = xg train.shape[1]-1
         xgb matrix = xgb.DMatrix(data=tr x, label=y)
         m1 = xgb.train(params=params, dtrain=xgb matrix)
         xt= xqb.DMatrix(data=x_test)
         preds = m1.predict(xt, output margin=False)
         for i in range(0, len(preds)):
             if preds[i] > 0.5:
                 preds[i]=1
             else:
                 preds[i]=0
         print evaluate(preds, y test)
         print preds
         0.509404254107
         [1. 1. 1. ..., 1. 1. 1.]
In [53]: def transform(preds):
             for i in range(0, len(preds)):
                 if preds[i] > 0.5:
                     preds[i]=1
                 else:
                     preds[i]=0
             return preds
         def cv(params, estimators):
             start = time.time()
             expected err = [0]*len(params);
             expected err stump = 0;
             xabs=[]
             for i in range(0, 10):
                 print i*10, '% complete'
                 x_train, x_test, y_train, y_test = sk.model_selection.train_t
         est_split(tr_x, y, test_size=test_size, random_state=seed)
                 x train = xgb.DMatrix(data=x train, label=y train)
                 x test = xgb.DMatrix(data=x test)
                 for i in range(0, len(params)):
                     xgbs.append(xgb.train(params[i], dtrain=x_train, num_boos
         t round=estimators[i]))
                     expected err[i] += evaluate(transform(xgbs[i].predict(x t
         est)), y test)
             for i in range(0, len(params)):
                 expected_err[i] = expected_err[i]/10
                 print 'Model ', i, ": ", expected_err[i]
             run time = time.time()-start
             print 'run time: ', int(run time)/60, 'm', run time%60, 's'
```

```
In [54]: #params = {'max_depth':5, 'learning_rate':0.3, 'n_estimators':1000,
    'objective':'binary:logistic', 'updater':'grow_gpu'}
    #param2 = {'max_depth':1, 'learning_rate':0.3, 'n_estimators':1000,
    'objective':'binary:logistic', 'updater':'grow_gpu'}
    #models = [params, param2]
    #estimators = [1000, 1000]
    #cv(models, estimators)
```

```
In [55]: #no gpu
#params = {'max_depth':5, 'learning_rate':0.3, 'n_estimators':1000,
    'objective':'binary:logistic'}
#param2 = {'max_depth':1, 'learning_rate':0.3, 'n_estimators':1000,
    'objective':'binary:logistic'}
#models = [params, param2]
#estimators = [1000, 1000]
#cv(models, estimators)
```

Apparently, GPU really slows down the models.

Stump is a little better, lets work with it.

Model 0: 0.5

Model 1: 0.516098481114 run time: 0: 12.8191699982 s

Samller learning rate and increased estimators both seem to do worse.

```
In [57]: xgb_s25 = xgb.XGBClassifier(max_depth=1, learning_rate=0.1, n_estimat
    ors=25)
    xgb_s75 = xgb.XGBClassifier(max_depth=1, learning_rate=0.75, n_estimators=50)
    cross_val([xgb_s25, xgb_s75])
```

Model 0: 0.5

Model 1: 0.534483506517 run time: 0: 8.12539982796 s

So increased learning rate leads to better results.

```
In [58]: xgb75 = xgb.XGBClassifier(max_depth = 3, learning_rate=0.75, n_estima
    tors =50)
    xgb75_25 = xgb.XGBClassifier(max_depth = 1, learning_rate=0.75, n_est
    imators=25)
    cross_val([xgb75, xgb75_25])
```

Model 0: 0.565325455195 Model 1: 0.535515157092 run time: 0: 8.99209594727 s

Increased max depth is really good. Let's try playing with the number of estimators.

```
In [59]: xgb100 = xgb.XGBClassifier(max_depth = 3, learning_rate=0.75, n_estim
    ators =100)
    xgb5 = xgb.XGBClassifier(max_depth = 5, learning_rate=0.75, n_estimat
    ors =50)
    cross_val([xgb100, xgb5])
```

Model 0: 0.572970492033 Model 1: 0.575680519133 run time: 0: 16.2646460533 s

What happens at learning rate 1?

```
In [60]: xgb1_3 = xgb.XGBClassifier(max_depth = 3, learning_rate=1, n_estimato
    rs =50)
    xgb1_5 = xgb.XGBClassifier(max_depth = 5, learning_rate=1, n_estimato
    rs =50)
    cross_val([xgb1_3, xgb1_5])
```

Model 0: 0.578128744909 Model 1: 0.593599779167 run time: 0: 11.307612896 s

Allowing a higher max depth just gave a huge increase in performance when coupled with a higher learning rate.

```
In [61]: xgb100 = xgb.XGBClassifier(max_depth = 5, learning_rate=1, n_estimato
    rs=100)
    xgb15 = xgb.XGBClassifier(max_depth=5, learning_rate=1.5, n_estimator
    s=50)
    cross_val([xgb100, xgb15])
```

Model 0: 0.626343347508 Model 1: 0.595139477668 run time: 0: 17.580124855 s

```
In [62]: xgb10 = xgb.XGBClassifier(max_depth = 10, learning_rate=1, n_estimato
    rs=100)
    xgb200 = xgb.XGBClassifier(max_depth = 5, learning_rate=1, n_estimato
    rs=200)
    cross_val([xgb10, xgb200])
```

Model 0: 0.638700150947 Model 1: 0.635585920128 run time: 0: 38.0626480579 s

Increasing estimators, and depth of estimators worked wonders!!! Eventually I will have to do more than guess at changes to the model...

```
In [64]: xgb10 = xgb.XGBClassifier(max_depth = 10, learning_rate=1, n_estimato
    rs=200)
    xgb200 = xgb.XGBClassifier(max_depth = 5, learning_rate=1, n_estimato
    rs=1000)
    cross_val([xgb10, xgb200])
```

```
KeyboardInterrupt
                                          Traceback (most recent call
last)
<ipython-input-64-9e458cf20d0e> in <module>()
      1 xgb10 = xgb.XGBClassifier(max depth = 10, learning rate=1, n
estimators=200)
      2 xgb200 = xgb.XGBClassifier(max depth = 5, learning rate=1, n
estimators=1000)
----> 3 cross val([xgb10, xgb200])
<ipython-input-51-f867f104bbd8> in cross val(xgbs)
                x train, x test, y train, y test =
sk.model selection.train test split(tr x, y, test size=test size, ran
dom state=seed)
                for i in range(0, len(xgbs)):
---> 10
                    xgbs[i].fit(x train, y train)
                                                     #stump
     11
                    expected err[i] += evaluate(xgbs[i].predict(x tes
t), y_test)
     12
            for i in range(0, len(xgbs)):
/home/justin/anaconda2/lib/python2.7/site-packages/xgboost-0.6-py2.7.
egg/xgboost/sklearn.pyc in fit(self, X, y, sample weight, eval set, e
val metric, early stopping rounds, verbose)
                                      early stopping rounds=early sto
pping_rounds,
    463
                                      evals result=evals result,
obj=obj, feval=feval,
--> 464
                                      verbose eval=verbose)
    465
    466
                self.objective = xgb_options["objective"]
/home/justin/anaconda2/lib/python2.7/site-packages/xgboost-0.6-py2.7.
egg/xgboost/training.pyc in train(params, dtrain, num boost round, ev
als, obj, feval, maximize, early stopping rounds, evals result, verbo
se eval, xgb model, callbacks, learning rates)
    202
                                   evals=evals,
    203
                                   obj=obj, feval=feval,
--> 204
                                   xgb model=xgb model, callbacks=cal
lbacks)
    205
    206
/home/justin/anaconda2/lib/python2.7/site-packages/xgboost-0.6-py2.7.
egg/xgboost/training.pyc in _train_internal(params, dtrain, num_boost
round, evals, obj, feval, xgb model, callbacks)
     72
                # Skip the first update if it is a recovery step.
     73
                if version % 2 == 0:
---> 74
                    bst.update(dtrain, i, obj)
     75
                    bst.save rabit checkpoint()
     76
                    version += 1
/home/justin/anaconda2/lib/python2.7/site-packages/xgboost-0.6-py2.7.
egg/xgboost/core.pyc in update(self, dtrain, iteration, fobj)
    814
                if not isinstance(dtrain, DMatrix):
                    raise TypeError('invalid training matrix: {}'.for
    815
mat(type(dtrain). name ))
```

```
self. validate_features(dtrain)
--> 816
    817
    818
                if fobj is None:
/home/justin/anaconda2/lib/python2.7/site-packages/xgboost-0.6-py2.7.
egg/xgboost/core.pyc in validate features(self, data)
   1176
                else:
   1177
                    # Booster can't accept data with different featur
e names
-> 1178
                    if self.feature names != data.feature names:
                        dat missing = set(self.feature names) - set(d
   1179
ata.feature names)
                        my missing = set(data.feature names) - set(se
   1180
lf.feature names)
KeyboardInterrupt:
```

Its incredible to me that these keep improving. But those last 2 models took at least 5 minutes to cross validate. Might as well go for gold though.

It felt like this took 10 minutes to run, which is terribly disapointing for a 1% increase in accuracy. I have gone back and patched in a runtime counter.

So accuracy seems to increase with max depth and n_estimators, though runtime gets really bad if we increase these with reckless abandon. I will try one more run with max_depth=15 and n_estimators = 1000, but I expect minimal (if any) increase over 10/1000.

69%!!!! This jumped 1.5%, which was much more than I expected. Why stop bruteforcing? (Besides the runtime...)

```
In [ ]: xgb15 1000.fit(tr x, y)
         final preds = xgb15 1000.predict(test x fin)
In [ ]: print final preds.shape
         print id.shape
         print test x fin.shape
         get submission(final preds, 'boosted preds')
In [57]: | lgr = {'booster':'gblinear'}
         estimators = [10, 50, 100, 500]
         cv([lgr, lgr, lgr, lgr], estimators)
         0 % complete
         10 % complete
         20 % complete
         30 % complete
         40 % complete
         50 % complete
         60 % complete
         70 % complete
         80 % complete
         90 % complete
         Model 0: 0.64467053563
         Model 1: 0.662043189551
         Model 2: 0.667301562264
         Model 3: 0.665461504251
         run time: 0 m 38.6918599606 s
```

Linear regressison models seem to be more accurate by default. Stacking estimators does not seem to help very much.

```
In [59]: | lgr = {'booster':'gblinear'}
         estimators = [5, 10, 15, 20, 25, 50]
         cv([lgr, lgr, lgr, lgr, lgr], estimators)
         0 % complete
         10 % complete
         20 % complete
         30 % complete
         40 % complete
         50 % complete
         60 % complete
         70 % complete
         80 % complete
         90 % complete
        Model 0: 0.635913007473
        Model 1: 0.64467053563
        Model 2: 0.651283921893
        Model 3: 0.6552681218
        Model 4: 0.657897308157
        Model 5: 0.662043189551
         run time: 0 m 9.53281593323 s
```

We will stick with 50, as it seems to be the best tradeoff between speed and score.

Now lets add some weights to the model.

```
In [61]: | lgr ={}
          lgr['booster'] = 'gblinear'
          lgr['lambda'] = 0.1
          cv([lgr], [50])
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.652254010818
          run time: 0 m 4.95634007454 s
          cv data = xgb.DMatrix(data=tr x, label=y)
In [87]:
          res = []
          lambdas = []
          for i in range(0, 10):
              lgr['lambda'] = float(i)/10
              lambdas.append(float(i)/10)
              res.append(xgb.cv(lgr, dtrain = cv data, num boost round = 50, me
          trics={'error'}))
In [88]: type(res[0])
Out[88]: pandas.core.frame.DataFrame
In [89]:
          res[0].head()
Out[89]:
            test-error-mean
                          test-error-std train-error-mean
                                                      train-error-std
            0.063414
                           0.001144
                                       0.022857
                                                      0.000265
            0.061186
                           0.000799
                                                      0.000413
                                       0.019805
          2 0.062223
                           0.000653
                                                      0.000078
                                       0.018600
          3 0.062407
                           0.000604
                                       0.017776
                                                      0.000238
```

0.017532

0.000336

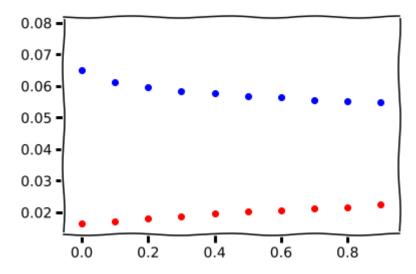
0.001142

0.062834

```
In [90]: testerr = []
    trainerr = []
    for df in res:
        testerr.append(df['test-error-mean'].mean(axis=0))
        trainerr.append(df['train-error-mean'].mean(axis=0))
```

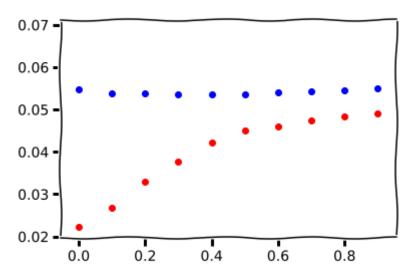
In [91]: **print** lambdaerr

```
In [92]: plt.scatter(lambdas, testerr, c='b')
   plt.scatter(lambdas, trainerr, c='r')
   plt.show()
```



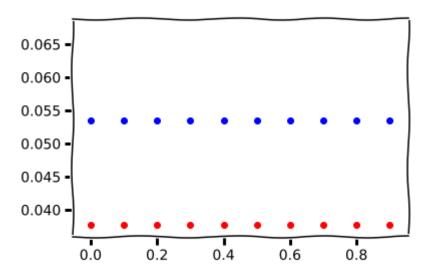
It seems lambda will be 0.9

```
In [95]:
         lgr['lambda'] = 0.9
         lgr['alpha'] = 0
         cv_data = xgb.DMatrix(data=tr_x, label=y)
         res = []
         alphas = []
         for i in range(0, 10):
             lgr['alpha'] = float(i)/10
             alphas.append(float(i)/10)
             res.append(xgb.cv(lgr, dtrain = cv data, num boost round = 50, me
         trics={'error'}))
         testerr = []
         trainerr = []
         for df in res:
             testerr.append(df['test-error-mean'].mean(axis=0))
             trainerr.append(df['train-error-mean'].mean(axis=0))
         plt.scatter(alphas, testerr, c='b')
         plt.scatter(alphas, trainerr, c='r')
         plt.show()
```



Test error remains constant, but it seems that training error increases with increase in alpha. I'll keep it at 0.3 for now.

```
In [97]: |gr['lambda'] = 0.9
         lgr['alpha'] = 0.3
         lgr['lambda bias'] = 0
         cv data = xgb.DMatrix(data=tr x, label=y)
         res = []
         lb = []
         for i in range(0, 10):
             lgr['lambda bias'] = float(i)/10
             lb.append(float(i)/10)
             res.append(xgb.cv(lgr, dtrain = cv data, num boost round = 50, me
         trics={'error'}))
         testerr = []
         trainerr = []
         for df in res:
             testerr.append(df['test-error-mean'].mean(axis=0))
             trainerr.append(df['train-error-mean'].mean(axis=0))
         plt.scatter(lb, testerr, c='b')
         plt.scatter(lb, trainerr, c='r')
         plt.show()
```



This will just stay at 0.

```
In [120]: class my model():
              def init (self, gbl params, gbt params, ngbl, ngbt):
                   self.lin = qbl params
                   self.tree = gbt params
              def fit(self, x, y):
                   data = xgb.DMatrix(data=x, label = y)
                   self.m1 = xgb.train(self.lin, dtrain = data, num boost round
          = 50)
                  preds = transform(self.ml.predict(xqb.DMatrix(data=x)))
                  err = []
                  for i in range(0, len(y)):
                       if y[i] != preds[i]:
                           err.append(1)
                       else:
                           err.append(0)
                   data = xgb.DMatrix(data=x, label=err)
                  self.m2 = xqb.train(self.tree, dtrain = data,
          num boost round=200)
              def predict(self, x):
                  data = xqb.DMatrix(data=x)
                   preds = transform(self.ml.predict(data))
                  err = transform(self.m2.predict(data))
                   preds = np.absolute(preds-err)
                   return preds
In [100]: | lgr
Out[100]: {'alpha': 0.3,
            'aplha': 5.0,
           'booster': 'gblinear',
           'lambda': 0.9,
           'lambda bias': 0.9}
In [102]: | lgr['lambda bias']=0
          cv([lgr], [50])
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.591536478016
          run time: 0 m 4.5983080864 s
```

```
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.609821383582
run time: 0 m 4.8096280098 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.609113095981
run time: 0 m 4.90501499176 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.608000604662
run time: 0 m 4.91266393661 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.605856462768
run time: 0 m 4.76365709305 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
```

```
90 % complete
Model 0: 0.600840612286
run time: 0 m 5.0142250061 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.59885815188
run time: 0 m 5.03225302696 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.596633169242
run time: 0 m 4.82712888718 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.596875691473
run time: 0 m 4.98250412941 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
60 % complete
70 % complete
80 % complete
90 % complete
Model 0: 0.594246505117
run time: 0 m 5.42937421799 s
0 % complete
10 % complete
20 % complete
30 % complete
40 % complete
50 % complete
```

```
60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.591536478016
          run time: 0 m 5.63348698616 s
In [105]: | gbl = {'booster':'gblinear'}
          cv([gbt], [50])
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.660607335257
          run time: 0 m 4.91333413124 s
In [123]: gbt = {'max depth':15, 'learing rate':1, 'objective':'binary:logisti
          cv([gbt, gbl], [200, 50])
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.632490968402
          Model 1: 0.661881508063
          run time: 0 m 42.7306771278 s
In [121]: lintree = my_model(gbl, gbt, 50, 200)
In [127]: cross_val([lintree])
          Model 0: 0.664395961906
          run time: 0 : 45.5462138653 s
```

Lol, it got worse...

```
In [128]:
          #some more gpu tests
          param = {'max_depth':3, 'learning_rate':0.3, 'objective':'binary:logi
          stic', 'updater':'grow gpu'}
          param2 = {'max depth':3, 'learning rate':0.3, 'objective':'binary:log
          istic'}
          cv([param], [50])
          cv([param2], [50])
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.507968399813
          run time: 0 m 19.4526128769 s
          0 % complete
          10 % complete
          20 % complete
          30 % complete
          40 % complete
          50 % complete
          60 % complete
          70 % complete
          80 % complete
          90 % complete
          Model 0: 0.507968399813
          run time: 0 m 6.60225105286 s
```

CPU still seems much faster than GPU

```
In [ ]:
```