In [2]:

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.svm import SVR
from sklearn.metrics import mean_squared_error
from sklearn.model_selection import RandomizedSearchCV, GridSearchCV
```

Lendo arquivos

In [3]:

0.1 SVM Regressor

0.2 Medida de erro

In [12]:

```
list_exp_C = range(-5, 15, 5)
list_exp_gamma = range(-15, 3, 3)
list epsilon = np.arange(0.05, 1, 0.5)
#Lista de parametros para utilizar nos demais
C list = []
gamma_list = []
best params = []
best score = 0
mse = 0
for exp gamma in list exp gamma:
    gamma = 2**exp gamma
    gamma_list.append(gamma)
    for exp C in list exp C:
        C = 2**exp C
        C list.append(C)
        for epilson in list epsilon:
            clf = SVR(gamma=gamma, C=C, epsilon=epilson)
            clf.fit(x train, y train)
            score = clf.score(x test,y test)
            if score > best score:
                best score = score
                best params = [gamma, C, epilson]
                mse = mean squared error(clf.predict(x test), y test)
print("Best Score: ",best_score ," with params: ", best_params," and MSE: ",mse)
C list = np.unique(C list)
gamma list = np.unique(gamma list)
list epsilon = np.unique(list epsilon)
```

Best Score: 0.7101346849881802 with params: [0.000244140625, 1024, 0.05] and MSE: 19.45487953760004

1 Random search

```
In [64]:
```

2 Grid seach

Model with rank: 3

Mean validation score: 0.723 (std: 0.040)

In [63]:

```
grid search = GridSearchCV(clf, param grid=param dist, cv=5)
grid search.fit(x train, y train)
report(grid search.cv results )
/home/pazeto/anaconda3/envs/MO431A/lib/python3.7/site-packages/sklear
n/model selection/ search.py:841: DeprecationWarning: The default of t
he `iid` parameter will change from True to False in version 0.22 and
will be removed in 0.24. This will change numeric results when test-se
t sizes are unequal.
  DeprecationWarning)
Model with rank: 1
Mean validation score: 0.746 (std: 0.040)
Parameters: {'C': 32768, 'epsilon': 1.0, 'gamma': 3.0517578125e-05}
Model with rank: 2
Mean validation score: 0.744 (std: 0.039)
Parameters: {'C': 32768, 'epsilon': 0.5, 'gamma': 3.0517578125e-05}
Model with rank: 3
Mean validation score: 0.744 (std: 0.034)
Parameters: {'C': 32768, 'epsilon': 0.1, 'gamma': 3.0517578125e-05}
```

Parameters: {'gamma': 3.0517578125e-05, 'epsilon': 1.0, 'C': 1024}

3 Otimização bayesiana

In [4]:

```
# !pip install hyperopt
from hyperopt import hp, tpe, fmin, space eval
def my svr fun(params):
    C, gamma, epsilon = params['C'], params['gamma'], params['epsilon']
    clf = SVR(gamma=gamma, C=C, epsilon=epsilon)
    clf.fit(x train, y train)
    score = clf.score(x test, y test)
    return -score
param_hyperopt= {
    'C': 2 ** hp.uniform('C', -5, 15),
    'gamma': 2 ** hp.uniform('gamma', -15, 3),
    'epsilon': hp.uniform('epsilon', 0.05, 1),
}
best_params = fmin(fn = my_svr_fun,
            space = param hyperopt, algo=tpe.suggest,
            \max \text{ evals} = 125)
print("Melhores paramtros: ", space eval(param hyperopt, best params))
print("Score: ", my_svr_fun(space_eval(param_hyperopt, best_params))*-1)
            | 125/125 [00:50<00:00, 2.48it/s, best loss: -0.825205
100%
8541104953]
Melhores paramtros: {'C': 18254.13344755772, 'epsilon': 0.15646708299
651826, 'gamma': 3.194344320388373e-05}
```

4 PSO

Score: 0.8252058541104953

In [160]:

```
from pyswarm import pso
#primeiro elemento é o mais baixo e mais alte de C, segundo de gamma e terceiro de
lb = [2**-5, 2**-15, 0.05]
ub = [2**15, 2**3, 1.0]
#dado que o pso vai esperar o menor valor e
# queremos o melhor score devemos negativar o score, o mais baixo terá melhores par
def my_svr_fun(params):
     print(params)
   C, gamma, epilson = params
    clf = SVR(gamma=gamma, C=C, epsilon=epilson)
    clf.fit(x_train, y_train)
    score = clf.score(x test,y test)
    return -score
best params, best score = pso(my svr fun, lb, ub, swarmsize=11, maxiter=11)
print("Parametros: ",best_params)
print("Score:", best_score*-1)
```

Stopping search: maximum iterations reached --> 11

Parametros: [1.86712162e+04 3.05175781e-05 5.29600172e-02]

Score: 0.8296045658080052

5 Simulated annealing

```
In [ ]:
```

```
from simanneal import Annealer
import random
class SearchBestParamsSVR(Annealer):
    def move(self):
        """permuta a lista de parametros possíveis"""
        random.shuffle(self.state['C'])
        random.shuffle(self.state['gamma'])
        random.shuffle(self.state['epsilon'])
    def energy(self):
        """Calcula o melhor score"""
        C, gamma, epilson = self.state['C'][0], self.state['gamma'][0], self.state[
        clf = SVR(gamma=gamma, C=C, epsilon=epilson)
        clf.fit(x train, y train)
        score = clf.score(x_test, y_test)
        return -score
param dist = {^{"}C": [2**-5, 2**0, 2**5, 2**10, 2**15],
              "gamma": [2**-15, 2**-10, 2**-5, 2**0, 2**5],
              "epsilon": [0.05, 0.1, 0.5, 0.8, 1.0]}
tsp = SearchBestParamsSVR(param dist)
params, score = tsp.anneal()
```

In [119]:

```
print("Parametros C gamma e epilson:",params['C'][0], params['gamma'][0], params['e
print("Score:", score*-1)
```

Pametros C gamma e epilson: 32768 3.0517578125e-05 0.05

Score: 0.8220660478366459

6 CMA-ES

84

5.1 88

```
# !pip install cma
import cma
# help(cma.fmin)
def my svr fun(params):
    C, gamma, epilson = params
    clf = SVR(gamma=gamma, C=C, epsilon=epilson)
    clf.fit(x_train, y_train)
    score = clf.score(x test,y test)
    return -score
lb = [2**-5, 2**-15, 0.05]
ub = [2**15, 2**3, 1]
opt = { 'bounds': [lb , ub], 'verb_disp':1000}
opt['scaling of variables'] = [2**10, 2**1, 0.01]
es = cma.fmin(my_svr_fun, [1, 1, 1], 2.5, opt)
print("Parametros C gamma e epilson: ", es[0])
print("Score: ", es[1]*-1)
(3 w,7)-aCMA-ES (mu w=2.3,w 1=58%) in dimension 3 (seed=769965, Wed Ma
y 1 11:49:26 2019)
Iterat #Fevals
                 function value axis ratio sigma min&max std t[m:
s1
           7 -1.838206063591308e-04 1.0e+00 2.08e+00
                                                       2e+00 2e+00 0:0
0.1
          14 1.480056132296959e-04 1.4e+00 1.93e+00
                                                      1e+00
                                                              2e+00 0:0
0.2
          21 -9.133457200733330e-03 1.8e+00 1.83e+00
                                                       1e + 00
                                                              2e+00 0:0
    3
0.3
         154 -6.326321821002134e-01 1.2e+01 1.54e-01
   22
                                                       1e-02
                                                              1e-01 0:0
3.5
         238 -6.475068819370783e-01 4.6e+01 7.24e-02
                                                       1e-03
                                                              6e-02 0:0
   34
7.6
         322 -7.111023370763685e-01 3.3e+02 1.39e-01
                                                       1e-03
                                                              2e-01 0:1
   46
3.1
         371 -7.681652760797661e-01 9.7e+02 7.36e-01
                                                       4e-03
                                                               2e+00 0:1
   53
9.7
         392 -7.774315533913859e-01 1.1e+03 9.40e-01
                                                       3e-03
                                                              2e+00 0:2
   56
7.1
         420 -7.824298371973549e-01 1.0e+03 8.63e-01
                                                       2e-03
                                                               1e+00 0:3
6.7
   64
         448 -7.975142264359517e-01 1.2e+03 6.91e-01
                                                       1e-03
                                                               1e+00 0:4
8.6
         469 -7.968387811299247e-01 1.3e+03 8.05e-01
                                                              2e+00 1:0
   67
                                                       9e-04
1.2
         490 -8.010360811367336e-01 1.8e+03 7.51e-01
   70
                                                       7e-04
                                                              1e+00 1:1
3.6
         518 -7.956126969487143e-01 1.6e+03 9.39e-01
                                                       1e-03
                                                               1e+00 1:2
9.4
   78
         546 -8.042409170371525e-01 1.6e+03 1.31e+00
                                                       2e-03
                                                               2e+00 1:4
6.8
         567 -7.969227474040904e-01 1.7e+03 1.28e+00
                                                       1e-03
                                                               1e+00 2:0
   81
4.8
```

588 -8.013909088537767e-01 2.0e+03 1.39e+00

616 -8.018597706470010e-01 1.5e+03 1.52e+00

1e-03

1e-03

2e+00 2:2

1e+00 2:4

```
466
        3262 -8.256966351184095e-01 2.6e+02 6.48e-02 3e-11 7e-05 53:
07.0
  473
        3311 -8.256993397236930e-01 2.3e+02 7.05e-02
                                                      3e-11
                                                            6e-05 54:
25.5
  482
        3374 -8.256917416376404e-01 1.1e+02 1.13e-01
                                                      7e-11
                                                             5e-05 55:
50.6
  490
        3430 -8.256894822402588e-01 9.9e+01 9.33e-02 4e-11
                                                            4e-05 57:
10.6
        3486 -8.257054469894023e-01 7.9e+01 1.11e-01
  498
                                                      5e-11 3e-05 58:
38.8
        3556 -8.257005453302735e-01 9.8e+01 1.53e-01
  508
                                                      5e-11
                                                             4e-05 60:
00.4
  519
        3633 -8.256941680231893e-01 2.3e+02 2.29e-01
                                                      6e-11
                                                            7e-05 61:
22.9
  530
       3710 -8.257020333945037e-01 2.1e+02 1.43e-01
                                                      3e-11
                                                            3e-05 62:
47.5
        3787 -8.256941719645042e-01 3.0e+02 2.00e-01
                                                      3e-11 7e-05 64:
  541
13.0
        3871 -8.257005250102156e-01 3.4e+02 2.49e-01
  553
                                                      4e-11
                                                             5e-05 65:
44.5
       3955 -8.257050357771927e-01 4.4e+02 1.10e-01 1e-11
                                                             2e-05 67:
  565
17.8
       4039 -8.257007381794392e-01 1.4e+03 4.05e-02 4e-12
  577
                                                            8e-06 68:
49.3
        4123 -8.256982696352093e-01 1.1e+03 4.68e-02
                                                      4e-12
                                                             7e-06 70:
  589
18.8
        4207 -8.256983777314512e-01 2.1e+03 6.43e-02
  601
                                                      7e-12
                                                             1e-05 71:
51.3
  613
        4291 -8.256978744085418e-01 3.4e+03 7.51e-02
                                                      8e-12
                                                             1e-05 73:
23.3
        4375 -8.256909369531588e-01 4.4e+03 6.49e-02
                                                      5e-12 8e-06 74:
  625
55.1
        4466 -8.257026310997099e-01 2.9e+03 1.03e-01 4e-12
  638
                                                             7e-06 76:
32.5
  650
        4550 -8.256973709038034e-01 4.3e+03 6.20e-02
                                                      2e-12 4e-06 78:
01.5
termination on tolstagnation=174 (Wed May 1 13:07:30 2019)
final/bestever\ f-value = -8.257023e-01\ -8.257230e-01
incumbent solution: [19397.06710389928, 3.051757888265911e-05, 0.36033
459459511381
std deviation: [0.0005189705279846922, 4.266018454044987e-12, 3.936943
727446785e-081
Parametros C gamma e epilson: [1.94095292e+04 3.05266127e-05 3.605850
58e-011
Score: 0.8257229759217513
```