## absorp\_ratio\_m2\_ex5

October 25, 2018

## 1 Course Project: Trading Strategy based on PCA

Welcome to your course project. This exercise gives you a hands-on experience to use PCA to:

- construct eigen-portfolios
- implement a measure of market systemic risk
- develop simple trading strategy

**Instructions:** - You will be using Python 3. - Avoid using for-loops and while-loops, unless you are explicitly told to do so. - Do not modify the (# GRADED FUNCTION [function name]) comment in some cells. Your work would not be graded if you change this. Each cell containing that comment should only contain one function. - After coding your function, run the cell right below it to check if your result is correct.

After this assignment you will: - Be able to use PCA to construct eigen-portfolios - Be able to use PCA to calculate a measure of market systemic risk - Be able to implement and analyze performance of portfolio strategy

Let's get started!

## 1.1 About iPython Notebooks

iPython Notebooks are interactive coding environments embedded in a webpage. You will be using iPython notebooks in this class. You only need to write code between the ### START CODE HERE ### and ### END CODE HERE ### comments. After writing your code, you can run the cell by either pressing "SHIFT"+"ENTER" or by clicking on "Run Cell" (denoted by a play symbol) in the upper bar of the notebook.

We will often specify "( X lines of code)" in the comments to tell you about how much code you need to write. It is just a rough estimate, so don't feel bad if your code is longer or shorter.

## **1.2 1 - Packages**

First, let's run the cell below to import all the packages that you will need during this assignment. - numpy is the fundamental package for scientific computing with Python. - pandas Python data analysis library - pandas scikit-learn - machine learning in Python. - matplotlib is a famous library to plot graphs in Python.

```
In [1]: import pandas as pd
     import numpy as np
```

```
import sklearn.decomposition
        import tensorflow as tf
        from tensorflow.contrib.layers import fully_connected
        import sys
        import matplotlib.pyplot as plt
        %matplotlib inline
        print("Package Versions:")
        print(" scikit-learn: %s" % sklearn.__version__)
        print(" tensorflow: %s" % tf.__version__)
        sys.path.append("..")
        import grading
        try:
            import sklearn.model_selection
            import sklearn.linear_model
        except:
            print("Looks like an older version of sklearn package")
        try:
            import pandas as pd
            print(" pandas: %s"% pd.__version__)
        except:
            print("Missing pandas package")
Package Versions:
  scikit-learn: 0.18.1
  tensorflow: 1.8.0
 pandas: 0.19.2
In [2]: ### ONLY FOR GRADING. DO NOT EDIT ###
        submissions=dict()
        assignment_key="LztgGBBtEeiaYgrsftMrjA"
        all_parts=["oZXnf", "ahjZa", "9tUbW", "wjLi0"]
        ### ONLY FOR GRADING. DO NOT EDIT ###
In [75]: # COURSERA_TOKEN = # the key provided to the Student under his/her email on submission
         # COURSERA_EMAIL = # the email
         COURSERA_TOKEN="Y3jMwrYGzVHMxG7K"
         COURSERA_EMAIL="cilsya@yahoo.com"
```

**Dataset: daily prices of stocks from S&P 500 index** For this exercise we will be working with S&P 500 Index stock prices dataset. The following cell computes returns based for a subset of S&P 500 index stocks. It starts with stocks price data:

```
In [4]: import os
        # load dataset
        asset_prices = pd.read_csv(os.getcwd() + '/data/spx_holdings_and_spx_closeprice.csv',
                             date_parser=lambda dt: pd.to_datetime(dt, format='\frac{\psi Y-\frac{\psi n}{\pm -\psi d}'}{\pm -\psi d}'),
                             index_col = 0).dropna()
        n_stocks_show = 12
        print('Asset prices shape', asset_prices.shape)
        asset_prices.iloc[:, :n_stocks_show].head()
Asset prices shape (3493, 419)
Out [4]:
                                  AΑ
                                        AAPL
                                                 ABC
                                                          ABT
                                                                  ADBE
                          Α
                                                                            ADI \
        2000-01-27 46.1112 78.9443 3.9286 4.5485
                                                     13.7898
                                                              15.6719 48.0313
        2000-01-28 45.8585 77.8245 3.6295 4.5485
                                                      14.2653
                                                               14.3906 47.7500
        2000-01-31 44.5952 78.0345 3.7054 4.3968
                                                      14.5730
                                                               13.7656
                                                                        46.7500
        2000-02-01 47.8377 80.7640 3.5804 4.5333
                                                      14.7128
                                                               13.9688
                                                                        49.0000
        2000-02-02 51.5434 83.4934 3.5290 4.5788
                                                     14.7968
                                                               15.3281
                                                                        48.1250
                        ADM
                                 ADP
                                        ADSK
                                                  AEE
                                                           AEP
        2000-01-27 10.8844 39.5477 8.1250 32.9375 33.5625
        2000-01-28 10.7143 38.5627 7.7188 32.3125 33.0000
        2000-01-31 10.6576 37.3807 7.6406 32.5625
                                                       33.5000
        2000-02-01 10.8844 37.9717 7.9219 32.5625
                                                       33.6875
        2000-02-02 10.6576 35.9032 7.9688 32.5625
                                                       33.6250
1.2.1 Calculate daily log-returns
In [5]: asset_returns = np.log(asset_prices) - np.log(asset_prices.shift(1))
        asset_returns = asset_prices.pct_change(periods=1)
        asset_returns = asset_returns.iloc[1:, :]
        asset_returns.iloc[:, :n_stocks_show].head()
Out[5]:
                                    AA
                                            AAPL
                                                       ABC
                                                                 ABT
                                                                          ADBE \
                           Α
        2000-01-28 -0.005480 -0.014185 -0.076134 0.000000 0.034482 -0.081758
        2000-01-31 -0.027548  0.002698  0.020912 -0.033352
                                                           0.021570 -0.043431
        2000-02-01 0.072710 0.034978 -0.033735 0.031045
                                                           0.009593 0.014761
        2000-02-02 0.077464 0.033795 -0.014356 0.010037
                                                            0.005709 0.097310
        2000-02-03 0.016340 -0.031014 0.045537 -0.006617 0.005670 0.126402
                         ADI
                                   ADM
                                             ADP
                                                      ADSK
                                                                 AEE
                                                                           AEP
        2000-01-28 -0.005857 -0.015628 -0.024907 -0.049994 -0.018975 -0.016760
        2000-01-31 -0.020942 -0.005292 -0.030651 -0.010131
                                                            0.007737 0.015152
        2000-02-01 0.048128 0.021281
                                       0.015810 0.036816
                                                           0.000000 0.005597
        2000-02-02 -0.017857 -0.020837 -0.054475 0.005920
                                                           0.000000 -0.001855
        2000-02-03 0.098701 0.000000 0.067217 0.035288 0.011516 0.033457
In [6]: def center_returns(r_df):
```

.....

Normalize, i.e. center and divide by standard deviation raw asset returns data

```
Arguments:
           r\_df -- a pandas.DataFrame of asset returns
           normed_df -- normalized returns
           mean_r = r_df.mean(axis=0)
           sd_r = r_df.std(axis=0)
           normed_df = (r_df - mean_r) / sd_r
           return normed_df
In [7]: normed_r = center_returns(asset_returns)
       normed_r.iloc[:, :n_stocks_show].head()
Out [7]:
                                                      ABC
                                                                ABT
                                   AA
                                           AAPL
                                                                         ADBE
        2000-01-28 -0.190054 -0.513710 -2.714709 -0.049779 2.182933 -2.684131
        2000-01-31 -0.898232 0.096888 0.688156 -1.757230 1.355644 -1.438899
       2000-02-01 2.319164 1.264327 -1.227995 1.539597 0.588289 0.451774
       2000-02-02 2.471738 1.221529 -0.548494 0.464060 0.339454 3.133764
        2000-02-03 0.510174 -1.122380 1.551619 -0.388563 0.336944 4.078966
```

Now we are ready to compute Absorption Ratio(AR). We do so by defining a moving look back window over which we collect returns for computing PCA. We start off from the earliest historical data and march forward moving by step\_size, which we also choose arbitrary. For each such window we compute PCA and AR, fixing in advance number of components in the enumerator. Specifically, for we use the following hyper-parameters:

ADP

2000-01-28 -0.212461 -0.766996 -1.540731 -1.803947 -1.372991 -0.994169 2000-01-31 -0.720771 -0.279098 -1.891884 -0.391433 0.547298 0.871919 2000-02-01 1.606541 0.975244 0.948126 1.272129 -0.008894 0.313197 2000-02-02 -0.616815 -1.012898 -3.348109 0.177340 -0.008894 -0.122594

ADSK

AEE

ADM

ADI

**Part 1 (Implement exponentially-weighted) Instructions:** Implement exponent\_weighting function which returns a sequence of  $w_i$  as np.array. See below:

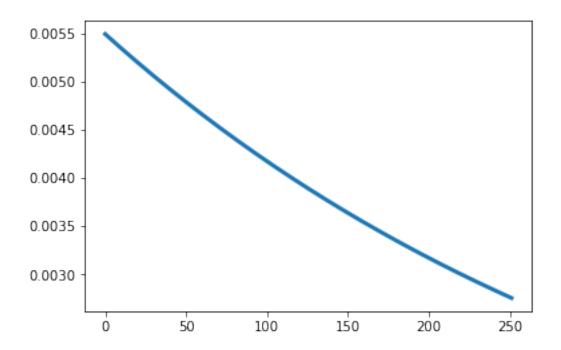
Define sequence of  $X_i$  where  $j \subset [N, 0]$ , an integer taking all values in the interval from 0 to N

$$X_j = e^{-\frac{log(2)}{H} \times j}$$

where H is half-life which determines the speed of decay, and log is natural log function Then a sequence of exponentially decaying weights  $w_i$  is defined as

$$w_j = \frac{X_j}{\sum\limits_{j=0}^{N} X_j}$$

```
In [13]: # GRADED FUNCTION: exponent_weighting# GRADE
         def exponent_weighting(n_periods, half_life = 252):
             Calculate exponentially smoothed normalized (in probability density function sense)
             Arguments:
             n_periods -- number of periods, an integer, N in the formula above
             half_life -- half-life, which determines the speed of decay, h in the formula
             Return:
             exp_probs -- exponentially smoothed weights, np.array
             11 11 11
             exp_probs = np.zeros(n_periods) # do your own calculation instead of dummy zero arr
             ### START CODE HERE ### ( 3 lines of code)
             ### ...
             for j in range(n_periods):
                 frac_temp = -1*(np.log(2)/half_life)*j
                 xj = np.exp(frac_temp)
                 exp\_probs[j] = xj
             denominator = exp_probs.sum()
             exp_probs = exp_probs / denominator
             return exp_probs
             ### END CODE HERE ###
In [14]: exp_probs = exponent_weighting(252*1)
         plt.plot(exp_probs, linewidth=3)
Out[14]: [<matplotlib.lines.Line2D at 0x7f8aa038cb70>]
```



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```
Out[15]: array([ 0.00549361, 0.00547852,
                                         0.00546347, 0.00544846,
                                                                  0.0054335 ,
                0.00541857, 0.00540369,
                                         0.00538885, 0.00537404,
                                                                 0.00535928,
                0.00534456, 0.00532988,
                                         0.00531524, 0.00530064,
                                                                 0.00528608,
                0.00527156, 0.00525708, 0.00524264, 0.00522824,
                                                                 0.00521388,
                0.00519956, 0.00518528,
                                         0.00517103, 0.00515683,
                                                                  0.00514267,
                0.00512854, 0.00511445,
                                         0.0051004 , 0.00508639,
                                                                  0.00507242,
                0.00505849, 0.00504459, 0.00503074, 0.00501692,
                                                                 0.00500314,
                0.0049894 , 0.00497569, 0.00496202, 0.00494839, 0.0049348 ,
                                        0.00489425, 0.00488081, 0.0048674,
                0.00492125, 0.00490773,
                0.00485403, 0.0048407,
                                         0.0048274 , 0.00481414,
                                                                 0.00480092,
                0.00478773, 0.00477458,
                                       0.00476146, 0.00474838, 0.00473534,
                0.00472233, 0.00470936, 0.00469643, 0.00468353, 0.00467066,
```

```
0.00465783, 0.00464504, 0.00463228, 0.00461956, 0.00460687,
                0.00459421, 0.00458159, 0.00456901, 0.00455646, 0.00454394,
                0.00453146, 0.00451901, 0.0045066, 0.00449422, 0.00448188,
                0.00446957, 0.00445729, 0.00444505, 0.00443284, 0.00442066,
                0.00440852, 0.00439641, 0.00438433, 0.00437229, 0.00436028,
                0.0043483, 0.00433636, 0.00432445, 0.00431257, 0.00430072,
                0.00428891, 0.00427713, 0.00426538, 0.00425367, 0.00424198,
                0.00423033, 0.00421871, 0.00420712, 0.00419557, 0.00418404])
In [16]: def absorption_ratio(explained_variance, n_components):
            Calculate absorption ratio via PCA. absorption_ratio() is NOT to be used with Auto-
            Arguments:
            explained_variance -- 1D np. array of explained variance by each pricincipal components
            n_components -- an integer, a number of principal components to compute absorption
            Return:
            ar -- absorption ratio
            ar = np.sum(explained_variance[:n_components]) / np.sum(explained_variance)
            return ar
```

**Part 2 (Implement Linear Auto-Encoder)** Linear Auto Encoder class has two fully connected layers and no activation functions between the layers.

**Instructions:** - fill missing code within LinearAutoEncoder class - in init() method of Linear-AutoEncoder setup neural network - **self.codings\_layer** is a fully connected layer with **n\_codings** neurons and no activation function - **self.outputs** is a fully connected layer with **n\_outputs** neurons and no activation function - define loss function as Mean Square Error between the outputs and inputs referenced by **self.X** in the code - use AdamOptimizer to optimize model parameters

```
# the inputs are n_inputs x n_inputs covariance matrices
    self.X = tf.placeholder(tf.float32, shape=[None, n_inputs, n_inputs])
    with tf.name_scope("lin_ae"):
        self.codings_layer = None
        self.outputs = None
        ### START CODE HERE ### ( 2 lines of code)
        self.codings_layer = tf.layers.dense(self.X, n_codings)
        self.outputs = tf.layers.dense(self.codings_layer, n_outputs)
        ### END CODE HERE ###
    with tf.name_scope("loss"):
        self.reconstruction_loss = None
        self.training_op = None
        ### START CODE HERE ### ( 4-5 lines of code)
        self.reconstruction_loss = tf.losses.mean_squared_error(self.X, self.output
        self.training_op = tf.train.AdamOptimizer(self.learning_rate).minimize(self
        ### END CODE HERE ###
        self.init = tf.global_variables_initializer()
def destroy(self):
    if hasattr(self, 'sess') and self.sess is not None:
        self.sess.close()
        self.sess = None
def absorption_ratio(self, test_input):
    Calculate absorption ratio based on already trained model
    if self.outputs is None:
        return test_input, 0.
    with self.sess.as_default(): # do not close session
        codings = self.codings_layer.eval(feed_dict={self.X: test_input})
        # calculate variance explained ratio
        result_ = self.outputs.eval(feed_dict={self.X: test_input})
        var_explained = np.sum(np.diag(result_.squeeze())) / np.sum(np.diag(test_in))
    return codings[0, :, :], var_explained
def next_batch(self, X_train, batch_size):
    \mathit{X\_train} - \mathit{np.array} of double of size \mathit{K} \ \mathit{x} \ \mathit{N} \ \mathit{x} \ \mathit{N}, where \mathit{N} is dimensionality of the
    batch_size - an integer, number of training examples to feed through the nwtwor
    y_batch = None
    selected_idx = np.random.choice(tuple(range(X_train.shape[0])), size=batch_size
```

```
X_batch = X_train[selected_idx, :, :]
    return X_batch, y_batch
def train(self, X_train, X_test, n_epochs=5, batch_size=2, verbose=False):
    train simple auto-encoder network
    :param X_train:
    :param X\_test:
    :param n_epochs: number of epochs to use for training the model
    :param batch_size:
    :return:
    11 11 11
    if self.outputs is None:
        return X_test, 0.
    n_examples = len(X_train) # number of training examples
    self.sess = tf.Session()
    # as_default context manager does not close the session when you exit the conte
    # and you must close the session explicitly.
    with self.sess.as_default():
        self.init.run()
        for epoch in range(n_epochs):
            n_batches = n_examples // min(n_examples, batch_size)
            for _ in range(n_batches):
                X_batch, y_batch = self.next_batch(X_train, batch_size)
                self.sess.run(self.training_op, feed_dict={self.X: X_batch})
            if verbose:
                # last covariance matrix from the training sample
                if X_train.shape[0] == 1:
                    mse_train = self.reconstruction_loss.eval(feed_dict={self.X: X_
                else:
                    mse_train = self.reconstruction_loss.eval(feed_dict={self.X: np
                mse_test = self.reconstruction_loss.eval(feed_dict={self.X: X_test})
                print('Epoch %d. MSE Train %.4f, MSE Test %.4f' % (epoch, mse_train
        # calculate variance explained ratio
        test_input = np.array([X_train[-1, :, :]])
        result_ = self.outputs.eval(feed_dict={self.X: test_input})
        var_explained = np.sum(np.diag(result_.squeeze())) / np.sum(np.diag(test_in))
        # print('Linear Auto-Encoder: variance explained: %.2f' % var_explained)
        codings = self.codings_layer.eval(feed_dict={self.X: X_test})
        # print('Done training linear auto-encoder')
    return codings[0, :, :], var_explained
```

```
In [24]: ### GRADED PART (DO NOT EDIT) ###
         ix offset = 1000
         stock_tickers = asset_returns.columns.values[:-1]
         assert 'SPX' not in stock_tickers, "By accident included SPX index"
         step_size = 60
         num\_samples = 5
         lookback\_window = 252 * 2 # in (days)
         num_assets = len(stock_tickers)
         cov_matricies = np.zeros((num_samples, num_assets, num_assets)) # hold training data
         # collect training and test data
         ik = 0
         for ix in range(ix_offset, min(ix_offset + num_samples * step_size, len(normed_r)), ste
             ret_frame = normed_r.iloc[ix_offset - lookback_window:ix_offset, :-1]
             print("time index and covariance matrix shape", ix, ret_frame.shape)
             cov_matricies[ik, :, :] = ret_frame.cov()
             ik += 1
         # the last covariance matrix determines the absorption ratio
         lin_ae = LinearAutoEncoder(n_inputs=num_assets, n_codings=200)
         np.array([cov_matricies[-1, :, :]]).shape
         lin_codings, test_absorp_ratio = lin_ae.train(cov_matricies[ : int((2/3)*num_samples),
                                                         np.array([cov_matricies[-1, :, :]]),
                                                         n_epochs=10,
                                                         batch_size=5)
         lin_codings, in_sample_absorp_ratio = lin_ae.absorption_ratio(np.array([cov_matricies[0
         ### GRADED PART (DO NOT EDIT) ###
time index and covariance matrix shape 1000 (504, 418)
time index and covariance matrix shape 1060 (504, 418)
time index and covariance matrix shape 1120 (504, 418)
time index and covariance matrix shape 1180 (504, 418)
time index and covariance matrix shape 1240 (504, 418)
In [25]: ### GRADED PART (DO NOT EDIT) ###
         part_2=[test_absorp_ratio, in_sample_absorp_ratio]
         try:
             part2 = " ".join(map(repr, part_2))
         except TypeError:
             part2 = repr(part_2)
         submissions[all_parts[1]]=part2
         grading.submit(COURSERA_EMAIL, COURSERA_TOKEN, assignment_key,all_parts[:2],all_parts,s
         [test_absorp_ratio, in_sample_absorp_ratio]
         ### GRADED PART (DO NOT EDIT) ###
```

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```
Out [25]: [0.39359208195496026, 0.39359208195496026]
In [26]: stock_tickers = asset_returns.columns.values[:-1]
         assert 'SPX' not in stock_tickers, "By accident included SPX index"
         half_life = 252
                                     # in (days)
         lookback_window = 252 * 2 # in (days)
         num_assets = len(stock_tickers)
                                # days : 5 - weekly, 21 - monthly, 63 - quarterly
         step_size = 1
         # require of that much variance to be explained. How many components are needed?
         var threshold = 0.8
         # fix 20% of principal components for absorption ratio calculation. How much variance (
         absorb\_comp = int((1 / 5) * num\_assets)
         print('Half-life = %d' % half_life)
         print('Lookback window = %d' % lookback_window)
         print('Step size = %d' % step_size)
         print('Variance Threshold = %d' % var_threshold)
         print('Number of stocks = %d' % num_assets)
         print('Number of principal components = %d' % absorb_comp)
Half-life = 252
Lookback window = 504
Step size = 1
Variance Threshold = 0
Number of stocks = 418
Number of principal components = 83
In [28]: # indexes date on which to compute PCA
         days_offset = 4 * 252
         num_days = 6 * 252 + days_offset
         pca_ts_index = normed_r.index[list(range(lookback_window + days_offset, min(num_days, l
         # allocate arrays for storing absorption ratio
         pca_components = np.array([np.nan]*len(pca_ts_index))
         absorp_ratio = np.array([np.nan]*len(pca_ts_index))
         lae_ar = np.array([np.nan]*len(pca_ts_index)) # absorption ratio computed by Auto-Enco
         # keep track of covariance matricies as we would need them for training Auto-Encoder
         buf_size = 5
         cov_matricies = np.zeros((buf_size, num_assets, num_assets))
         exp_probs = exponent_weighting(lookback_window, half_life)
         assert 'SPX' not in normed_r.iloc[:lookback_window, :-1].columns.values, "By accident i
```

**Instructions:** - on each loop iteration: - fit PCA to **cov\_mat** - use fitted pca model to pass values to absorption\_ratio(). The result of absorption ratio calculation goes into **absorp\_ratio** - compute

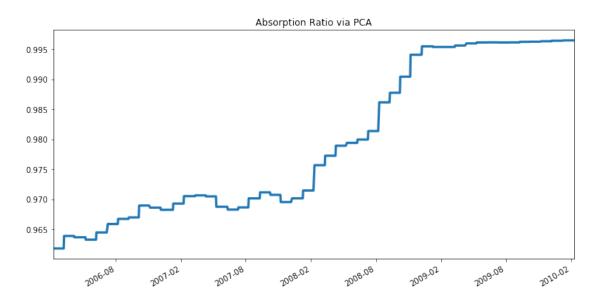
number of principal components it takes to explain at least **var\_threshold** of variance. The result of this calculation goes into **pca\_components** 

```
In [29]: # run the main loop computing PCA and absorption at each step using moving window of re
                        # run this loop using both exponentially weighted returns and equally weighted returns
                        import time
                        ik = 0
                        use_ewm = False
                        lin_ae = None
                        time_start = time.time()
                        for ix in range(lookback_window + days_offset, min(num_days, len(normed_r)), step_size)
                                   ret_frame = normed_r.iloc[ix - lookback_window:ix, :-1] # fixed window
                                   # ret_frame = normed_r.iloc[:ix, :-1] # ever-growing window
                                   if use_ewm:
                                              ret_frame = (ret_frame.T * exp_probs).T
                                  cov_mat = ret_frame.cov()
                                   circ idx = ik % buf size
                                  cov_matricies[circ_idx, :, :] = cov_mat.values
                                  if ik == 0 or ik % 21 == 0:
                                              ### START CODE HERE ### ( 4-5 lines of code)
                                              ### fit PCA, compute absorption ratio by calling absorption_ratio()
                                              ### store result into pca_components for grading
                                              \# https://www.coursera.org/learn/fundamentals-machine-learning-in-finance/discussions for the state of th
                                              \#\ http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.org/stable/modules/generated/sklearn.decomposition.PCA.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.http://scikit-learn.decomposition.pca.html
                                              pca=sklearn.decomposition.PCA().fit(cov_mat)
                                              absorp_ratio[ik] = absorption_ratio(pca.explained_variance_, absorb_comp)
                                              ### END CODE HERE ###
                                   else:
                                              absorp_ratio[ik] = absorp_ratio[ik-1]
                                              pca_components[ik] = pca_components[ik-1]
                                   if ik == 0 or ik % 252 == 0:
                                              if lin_ae is not None:
                                                        lin_ae.destroy()
                                              print('Trainging AE', normed_r.index[ix])
                                              lin_ae = LinearAutoEncoder(cov_mat.shape[0], absorb_comp)
                                              lin_codings, lae_ar[ik] = lin_ae.train(cov_matricies[:circ_idx + 1, :, :],
                                                                                                                                               np.array([cov_mat.values]),
                                                                                                                                               batch_size=2)
                                   else:
                                              lin_codings, lae_ar[ik] = lin_ae.absorption_ratio(np.array([cov_mat.values]))
```

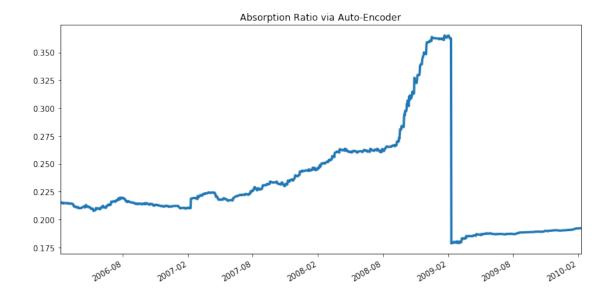
```
ik += 1
```

```
print ('Absorption Ratio done! Time elapsed: {} seconds'.format(time.time() - time_star
ts_pca_components = pd.Series(pca_components, index=pca_ts_index)
ts_absorb_ratio = pd.Series(absorp_ratio, index=pca_ts_index)
ts_lae_absorb_ratio = pd.Series(lae_ar, index=pca_ts_index)
```

Trainging AE 2006-02-07 00:00:00
Trainging AE 2007-02-09 00:00:00
Trainging AE 2008-02-11 00:00:00
Trainging AE 2009-02-10 00:00:00
Absorption Ratio done! Time elapsed: 35.51445770263672 seconds



In [31]: ts\_lae\_absorb\_ratio.plot(figsize=(12,6), title='Absorption Ratio via Auto-Encoder', lin
Out[31]: <matplotlib.axes.\_subplots.AxesSubplot at 0x7f8a2d360550>



Having computed daily (this means the step size is 1) Absorption Ratio times series, we further follow M. Kritzman to make use of AR to define yet another measure: AR Delta. In particular:

$$AR\delta = \frac{AR_{15d} - AR_{1y}}{AR\sigma_{1y}}$$

We use  $AR\delta$  to build simple portfolio trading strategy