model_per_device_and_time

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In [1]: import pandas as pd
        import numpy as np
        import calendar
        from bokeh.charts import output_notebook, Scatter, Bar, show, output_file, Line, BoxPlot
        from bokeh.plotting import figure
        from bokeh.layouts import row, column, gridplot
        from bokeh.charts import output_notebook, Scatter, Bar, show, output_file, Line, BoxPlot
        from bokeh.io import hplot
        from bokeh.models.ranges import Range1d
        from ML import filter_devices, build_deriv, resample_per_device, subsample_negatives
        from fft import fft_peak
        from sklearn.model_selection import cross_val_score
        from sklearn.ensemble import GradientBoostingClassifier, RandomForestClassifier
        from sklearn.pipeline import Pipeline
        from sklearn.decomposition import PCA
        from sklearn.preprocessing import Normalizer
        from sklearn.metrics import roc_curve, auc
        from sklearn.svm import SVC, NuSVC
        from sklearn.model_selection import LeavePGroupsOut, GroupShuffleSplit
        from sklearn.model_selection import train_test_split
        from sklearn.metrics import roc_curve, accuracy_score,precision_recall_curve, auc
        from sklearn.model_selection import GridSearchCV
        from sklearn.model_selection import StratifiedKFold
        output_notebook()
In [2]: INPUT="data/train.csv"
        dataset = pd.read_csv(INPUT,index_col=[0,1],parse_dates=[0])
0.1 Build Training set
In [3]: def pre_filter(df):
            res = df.copy()
            del res["attribute1"]
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del res["attribute3"]
            #del res["attribute5"]
            dt_list = ["attribute2"] #, "attribute3"]
            for c in dt_list:
                deriv = build_deriv(res,c)
                res["dt_%s" % c] = deriv
                res["dt2_%s" % c] = build_deriv(res,c,2)
            return res.fillna(0)
        def post_filter(df):
            res = df.copy()
            res = filter_devices(res)
            for col in res.columns:
                if "min" in col:
                    del res[col]
                if "std" in col:
                    del res[col]
            return res
In [4]: pre_dataset = pre_filter(dataset)
        #print feature_set.columns
        features = [f for f in pre_dataset.columns if "att" in f]
        def f_to_dict(feature):
            indexes = dict( ("avg_over%i_%s" % (i,feature), lambda df: df[:(i+1)].mean()) for i
            d = {
                    "min_%s" % feature:np.min,
                    "max_%s" % feature:np.max,
                    "mean_%s" % feature :np.mean,
                    "std_%s" % feature:np.std
            d.update(indexes)
            dft_list = ["attribute4", "attribute5", "attribute6", "attribute7", "attribute9"]
            if feature in dft_list:
                d["dft_p0_ind%s" % feature] = lambda r : fft_peak(r,p=0,index_no_value=True)
                d["dft_p0_val%s" % feature] = lambda r : fft_peak(r,p=0,index_no_value=False)
            return d
        agg_dict = dict( (f,f_to_dict(f)) for f in features )
In [5]: # bugfix: rolling aggregation after group by does not handle multiple aggregation per co
        # we fix this by flattening the aggregation dict and repeating the data within the data;
        final_columns = [k for c in agg_dict for k in agg_dict[c]]
        input_columns = [c for c in agg_dict for k in agg_dict[c]]
        flat_agg_dict = dict( (k,agg_dict[c][k]) for c in agg_dict for k in agg_dict[c])
        dup_dataset = pre_dataset[input_columns].sort_index(level="date").sort_index(level="devi
        dup_dataset.columns = final_columns
In [6]: #
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# instead of simply grouping by, we roll over the dataset...
        # using the "hacky" flat version, to avoid issues.
        # The hack increase the memory consumption quite a lot, but it should still be better the
        # explicitely building the windowed lines before aggregating over it.
        feature_set = resample_per_device(dup_dataset) \
            .groupby(level="device",as_index=False) \
            .rolling(window=360,min_periods=1) \
            .agg(flat_agg_dict) \
            .reset_index(level=0,drop=True)
        feature_set = post_filter(feature_set).sortlevel(level="device")
In [7]: #
        # feature filtering : removing features with a weak contribution to the last computed mo
        feat_filtering_thres = 5e-3
        try:
            kept_features = feature_imp[feature_imp.importance > feat_filtering_thres]
            print "threshold: %g, kept: %i features" % (feat_filtering_thres,kept_features.size
            filtered = feature_set.filter(items=feature_imp.index.sort_values())
            print "filtering devices"
            feature_set = filtered
        except:
            print "no feature filtering"
no feature filtering
In [8]: #
        # use label_window to expand label to neighboring days.
        \# basically, a mainrtenance x days before failure is still OK
        label_window = 7
        label_set = resample_per_device(dataset[["failure"]]) \
            .sortlevel(level="date",ascending=False) \
            .groupby(level="device",as_index=False) \
            .rolling(window=label_window, min_periods=1) \
            .reset_index(level=0,drop=True)
        label_set = filter_devices(label_set).sortlevel(level="device")
In [36]: #
         # subsampling the negatives, to balance classes
         negative_subsampling_fraction = 5e-2
         sub_label_set , sub_feature_set = subsample_negatives(negative_subsampling_fraction,lab
```

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126681 with 630 positives new size 6933
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try:

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In [9]: # compute groups (devices), to be used when splitting the training set (train/test)
        # This is useful to avoid the bias selection generated by a temporal model
        # (Basically, a device used in train cannot also be in test, because its attributes will
        # this is why the cross-val strategy must split by device, and not merely at random
        devices = sub_label_set.index.get_level_values("device")
        device_index = dict((device,i) for i,device in enumerate(devices.unique()))
        n_dev = len(device_index)
        device_groups = np.array(devices.to_series().map(device_index).tolist())
        NameError
                                                  Traceback (most recent call last)
        <ipython-input-9-35847931f59d> in <module>()
          3 # (Basically, a device used in train cannot also be in test, because its attributes
          4 # this is why the cross-val strategy must split by device, and not merely at random
    ----> 5 devices = sub_label_set.index.get_level_values("device")
          6 device_index = dict((device,i) for i,device in enumerate(devices.unique()))
          7 n_dev = len(device_index)
        NameError: name 'sub_label_set' is not defined
0.2 Run model
In [ ]: feature_mat = sub_feature_set.as_matrix()
        label_mat = sub_label_set.as_matrix().ravel()
        splitting_strategy = GroupShuffleSplit(n_splits=4,test_size=0.33)
        pca = PCA()
        norm = Normalizer()
        #model = Gradient BoostingClassifier()
        model = RandomForestClassifier()
        #model = SVC(probability=True)
        #pipeline= Pipeline([('normalize', norm),('reduce_dim', pca),("model",model)])
        pipeline= Pipeline([("model", model)])
        #pipeline = pipeline_optimizer._fitted_pipeline
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# use best parameters if available
            pipeline.set_params(**grid_result.best_params_)
            print "using last optimized model"
        except:
            print "no optim result, or bad ones: let's keep the default ones"
In [ ]: scores = cross_val_score(
            pipeline,
            feature_mat,
            label_mat,
            cv=splitting_strategy,
            groups = device_groups,
            verbose=1,
            scoring="f1",
            n_jobs=6)
        print "accurracy: %g, std(%g))" % (scores.mean(), scores.std())
0.2.1 Eval Model
In [15]: #
         # build the curve of PR, AUC on test and on train
         #
         X_train, X_test, Y_train, Y_test = train_test_split(feature_mat,label_mat,test_size=0.3
         # calculate the fpr and tpr for all thresholds of the classification
         fitted = pipeline.fit(X_train,Y_train)
         probs = fitted.predict_proba(X_test)
         preds = probs[:,1]
         preds_train = fitted.predict_proba(X_train)[:,1]
         fpr, tpr, threshold = roc_curve(Y_test, preds)
         fpr_train, tpr_train, threshold_train = roc_curve(Y_train, preds_train)
         roc_auc = auc(fpr, tpr)
         roc_auc_train = auc(fpr_train, tpr_train)
         precision, recall, ths = precision_recall_curve(Y_test, preds)
         precision_train, recall_train, ths_train = precision_recall_curve(Y_train, preds_train)
In [16]: #print "auc: %.2g, on train: %.2g" %(roc_auc, roc_auc_train)
         roc_df = pd.DataFrame({"fpr":fpr,"tpr":tpr}).set_index("fpr")
         pr_df = pd.DataFrame({"precision": precision, "recall":recall}).set_index("recall")
         roc_df["diag"] = roc_df.index
         pr_df["random"] = pr_df.precision.iloc[0]
         # roc curve
         roc_f = figure(width=400,height=400,title="roc, auc: %.2g, on train: %.2g" %(roc_auc,
         roc_f.xaxis.axis_label = "tpr"
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auc_range= Range1d(0,1)
roc_f.x_range = auc_range
roc_f.y_range = auc_range
roc_f.yaxis.axis_label = "fpr"
roc_f.cross(fpr,tpr,size=5)
roc_f.line(fpr,tpr,legend="roc")
roc_f.circle(fpr_train,tpr_train,size=5,color="red", line_width=1)
roc_f.line(fpr_train,tpr_train,color="red",legend="roc on train")
roc_f.line([0,1],[0,1], color="grey")
# pr curve
pr_f = figure(width=400,height=400,title="PR curve")
pr_f.xaxis.axis_label = "recall"
pr_f.yaxis.axis_label = "precision"
pr_f.cross(recall,precision,size=5)
pr_f.line(recall, precision, legend="PR")
pr_f.circle(recall_train,precision_train,size=5,color="red", line_width=1)
pr_f.line(recall_train,precision_train,color="red",legend="PR on train")
show(row(
    pr_f,
    roc_f
))
```

0.2.2 Debug: feature Importance

0.2.3 Hyperparameter optimisation: Grid search optim

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In [46]: # model : GradientBoostingClassifier, parameters:
         #loss : {deviance, exponential},
         #learning_rate : float, optional (default=0.1)
         \#n_estimators : int (default=100)
         #max_depth : integer, optional (default=3)
         #min_samples_split : int, float, optional (default=2)
         grids=dict()
         XDB_param_grid = {
             "model__loss": ["deviance", 'exponential'],
             "model__learning_rate" : [0.1],
             "model__n_estimators" : [100, 150],
             "model__max_depth" : [2,3,5],
             "model__min_samples_split" : [5,10,15]
         grids[GradientBoostingClassifier] = XDB_param_grid
In [47]: # C : penalty
         # kernel : linear, poly, rbf, sigmoid, precomputed
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SVC_param_grid = {
            'model__C': [1e-7,1e-6,1e-5,0.1],
             "model__kernel": ["rbf","linear"],
             #"model_degree" : [1,3,5], # polynomial degrees
             "model__gamma" : ["auto"], # kernel coef (rbf)
             #"coef0" # for poly, signmoid
             "model__tol" : [1e-3,1e-2]
         grids[SVC] = SVC_param_grid
In [48]: # model : RandomForestClassifier, parameters:
         # n_estimators : int (default=100)
         # criterion : "qini", "entropy"
         # max_features : auto , fraction
         # max_depth : integer, optional (default=3)
         # min_samples_split : int, float, optional (default=2)
         RF_param_grid = {
             "model__criterion": ["entropy"],
             "model__n_estimators" : [50,100,150],
             "model__max_features" : ["auto"],
             "model__max_depth" : [20,30,40],
             "model__min_samples_split" : [5,10,15]
         }
         grids[RandomForestClassifier] = RF_param_grid
In [49]: m = type(dict(pipeline.steps)["model"])
         param_grid=grids[m]
         splits = splitting_strategy.split(feature_mat,label_mat,device_groups)
         #kfold = StratifiedKFold(n_splits=6, shuffle=True)
         grid_search = GridSearchCV(
             pipeline,
             param_grid,
             scoring="f1",
             n_jobs=-1,
             verbose=1,
             cv=splits
         grid_result = grid_search.fit(feature_mat,label_mat)
         # summarize results
         print("Best: %f using %s" % (grid_result.best_score_, grid_result.best_params_))
         means = grid_result.cv_results_['mean_test_score']
         stds = grid_result.cv_results_['std_test_score']
         params = grid_result.cv_results_['params']
Fitting 4 folds for each of 27 candidates, totalling 108 fits
```