### Missing data in supervised ML

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https://github.com/brown-ccv/ODSC West 2019 (https://github.com/brown-ccv/ODSC West 2019)

#### About me

- · Born and raised in Hungary
- · Astrophysics PhD at MPIA, Heidelberg, Germany
- Postdoctoral researcher at MIT (still in astrophysics at the time)
- Started at Brown in December 2015 as a Data Scientist
- Promoted to Lead Data Scientist in 2017
- · Adjunct Lecturer in Data Science this semester
  - Teaching the course DATA1030: Hands-on data science to the DS master students at Brown

https://github.com/brown-ccv/ODSC\_West\_2019 (https://github.com/brown-ccv/ODSC\_West\_2019)

#### **Data Science at Brown**

- Center for Computation and Visualization
- · Institutional Data group
  - Data-driven decision support and predictive modeling for Brown's administrative units
  - Academic research on data-intensive projects
- OPEN POSITION more on this later

https://github.com/brown-ccv/ODSC\_West\_2019 (https://github.com/brown-ccv/ODSC\_West\_2019)

By the end of this workshop, you will be able to

- Describe the three main types of missingness patterns
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# Before we start, a few words on our dataset: kaggle house price

- · good for educational purposes
  - messy data that requires quite a bit of preprocessing
  - a nice mixture of continuous, ordinal, and categorical features, each feature type has missing values
- · lots of excellent kernels on kaggle
  - check them out here (https://www.kaggle.com/c/house-prices-advanced-regression-techniques)
- dataset and description available in repo
  - let's take a look!

### Missing values often occur in datasets

- survey data: not everyone answers all the questions
- medical data: not all tests/treatments/etc are performed on all patients
- sensor can be offline or malfunctioning
- customer data: not every user uses all features of an app

### Missing values are an issue for multiple reasons

#### Concenptual reason

- · missing values can introduce biases
  - bias: the samples (the data points) are not representative of the underlying distribution/population
  - any conclusion drawn from a biased dataset is also biased.
  - rich people tend to not fill out survey questions about their salaries and the mean salary estimated from survey data tend to be lower than true value

#### Practical reason

- missing values (NaN, NA, inf) are incompatible with sklearn
  - all values in an array need to be numerical otherwise sklearn will throw a ValueError
- there are a few supervised ML techniques that work with missing values (e.g., XGBoost, CatBoost)
  - we will cover those later today

### **Learning Objectives**

By the end of this workshop, you will be able to

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# Missing data patterns

- MCAR Missing Complete At Random
  - some people skip some survey questions by accident
- MAR Missing At Random
  - males are less likely to fill out a survey on depression
  - this has nothing to do with their level of depression after accounting for maleness
- MNAR Missing Not At Random
  - depressed people are less likely to fill out a survey on depression due to their level of depression

### **MCAR** test

- MCAR can be diagnosed with a statistical test (<u>Little, 1988</u> (<a href="https://www.tandfonline.com/doi/abs/10.1080/01621459.1988.10478722">https://www.tandfonline.com/doi/abs/10.1080/01621459.1988.10478722</a>))
  - python implementation available in the <u>pymice (https://github.com/RianneSchouten/pymice)</u> package or in the skipped slide

```
In [1]: # from the pymice package
        # https://github.com/RianneSchouten/pymice
        import numpy as np
        import pandas as pd
        import math as ma
        import scipy.stats as st
        def checks_input_mcar_tests(data):
            """ Checks whether the input parameter of class McarTests is correct
                    Parameters
                    data:
                        The input of McarTests specified as 'data'
                    Returns
                    _____
                    bool
                        True if input is correct
            if not isinstance(data, pd.DataFrame):
                print("Error: Data should be a Pandas DataFrame")
                return False
            if not any(data.dtypes.values == np.float):
                if not any(data.dtypes.values == np.int):
                    print("Error: Dataset cannot contain other value types than
         floats and/or integers")
                    return False
            if not data.isnull().values.any():
                print("Error: No NaN's in given data")
                return False
            return True
        def mcar test(data):
            """ Implementation of Little's MCAR test
            Parameters
             _____
            data: Pandas DataFrame
                An incomplete dataset with samples as index and variables as col
        umns
            Returns
            _____
            p value: Float
                This value is the outcome of a chi-square statistical test, test
        ing whether the null hypothesis
                 'the missingness mechanism of the incomplete dataset is MCAR' ca
        n be rejected.
            if not checks_input_mcar_tests(data):
                raise Exception("Input not correct")
```

```
dataset = data.copy()
    vars = dataset.dtypes.index.values
    n_var = dataset.shape[1]
    # mean and covariance estimates
    # ideally, this is done with a maximum likelihood estimator
    gmean = dataset.mean()
    gcov = dataset.cov()
    # set up missing data patterns
    r = 1 * dataset.isnull()
    mdp = np.dot(r, list(map(lambda x: ma.pow(2, x), range(n_var))))
    sorted mdp = sorted(np.unique(mdp))
    n pat = len(sorted mdp)
    correct mdp = list(map(lambda x: sorted mdp.index(x), mdp))
    dataset['mdp'] = pd.Series(correct_mdp, index=dataset.index)
    # calculate statistic and df
    pj = 0
    d2 = 0
    for i in range(n pat):
        dataset_temp = dataset.loc[dataset['mdp'] == i, vars]
        select_vars = ~dataset_temp.isnull().any()
        pj += np.sum(select_vars)
        select_vars = vars[select_vars]
        means = dataset temp[select vars].mean() - gmean[select vars]
        select_cov = gcov.loc[select_vars, select_vars]
        mj = len(dataset temp)
        parta = np.dot(means.T, np.linalg.solve(select cov, np.identity(
select cov.shape[1])))
        d2 += mj * (np.dot(parta, means))
    df = pj - n var
    # perform test and save output
    p value = 1 - st.chi2.cdf(d2, df)
    return p value
```

### **Takeaway**

- it can be challenging to infer the missingness pattern from an incomplete dataset
  - There is a statistical test to differentiate MCAR and MAR
  - MNAR is difficult/impossible to diagnose to the best of my knowledge
- · multiple patterns can be present in the data
  - even worse, multiple patterns can be present in one feature!
  - missing values in a feature can occur due to a mix of MCAR, MAR, MNAR

By the end of this workshop, you will be able to

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### Simple approaches for handling missing values

- 1) categorical/ordinal features: treat missing values as another category
  - missing values in categorical/ordinal features are not a big deal
- 2) continuous features: this is the tough part
  - sklearn's SimpleImputer
- 3) exclude points or features with missing values
  - might be OK

#### 1a) Missing values in a categorical feature

- YAY this is not an issue at all!
- Categorical feature needs to be one-hot encoded anyway
- Just replace the missing values with 'NA' or 'missing' and treat it as a separate category

### 1b) Missing values in a ordinal feature

- · this can be a bit trickier but usually fine
- · Ordinal encoder is applied to ordinal features
  - where does 'NA' or 'missing' fit into the order of the categories?
  - usually first or last
- if you can figure this out, you are golden

```
In [2]: # read the data
        import pandas as pd
        import numpy as np
        from sklearn.model_selection import train_test_split
        # Let's load the data
        df = pd.read csv('data/train.csv')
        # drop the ID
        df.drop(columns=['Id'],inplace=True)
        # the target variable
        y = df['SalePrice']
        df.drop(columns=['SalePrice'],inplace=True)
        # the unprocessed feature matrix
        X = df.values
        print(X.shape)
        # the feature names
        ftrs = df.columns
        (1460, 79)
In [3]: # let's split to train, test, and holdout
        X_other, X_holdout, y_other, y_holdout = train_test_split(df, y, test_si
        ze=0.2, random_state=0)
        X train, X test, y train, y test = train test split(X other, y other, te
        st size=0.25, random state=0)
        print(X train.shape)
        print(X test.shape)
        print(X holdout.shape)
        (876, 79)
        (292, 79)
```

(292, 79)

In [4]: # collect the various features cat ftrs = ['MSZoning','Street','Alley','LandContour','LotConfig','Neigh borhood', 'Condition1', 'Condition2', \ 'BldgType','HouseStyle','RoofStyle','RoofMatl','Exterior1st' , 'Exterior2nd', 'MasVnrType', 'Foundation', \ 'Heating','CentralAir','Electrical','GarageType','PavedDrive' , 'MiscFeature', 'SaleType', 'SaleCondition'] ordinal ftrs = ['LotShape','Utilities','LandSlope','ExterQual','ExterCon d','BsmtQual','BsmtCond','BsmtExposure',\ 'BsmtFinType1', 'BsmtFinType2', 'HeatingQC', 'KitchenQual', 'Functional', 'FireplaceQu', 'GarageFinish', \ 'GarageQual', 'GarageCond', 'PoolQC', 'Fence'] ordinal\_cats = [['Reg','IR1','IR2','IR3'],['AllPub','NoSewr','NoSeWa','E LO'],['Gtl','Mod','Sev'],\ ['Po','Fa','TA','Gd','Ex'],['Po','Fa','TA','Gd','Ex'],['N A', 'Po', 'Fa', 'TA', 'Gd', 'Ex'], \ ['NA','Po','Fa','TA','Gd','Ex'],['NA','No','Mn','Av','Gd' ],['NA','Unf','LwQ','Rec','BLQ','ALQ','GLQ'],\ ['NA','Unf','LwQ','Rec','BLQ','ALQ','GLQ'],['Po','Fa','T A','Gd','Ex'],['Po','Fa','TA','Gd','Ex'],\ ['Sal','Sev','Maj2','Maj1','Mod','Min2','Min1','Typ'],['N A','Po','Fa','TA','Gd','Ex'],\ ['NA', 'Unf', 'RFn', 'Fin'], ['NA', 'Po', 'Fa', 'TA', 'Gd', 'Ex'], ['NA','Po','Fa','TA','Gd','Ex'], ['NA', 'Fa', 'TA', 'Gd', 'Ex'], ['NA', 'MnWw', 'GdWo', 'MnPrv', 'G dPrv']] num ftrs = ['MSSubClass','LotFrontage','LotArea','OverallQual','OverallC ond','YearBuilt','YearRemodAdd',\ 'MasVnrArea', 'BsmtFinSF1', 'BsmtFinSF2', 'BsmtUnfSF', 'TotalBs mtSF', '1stFlrSF', '2ndFlrSF', \ 'LowQualFinSF','GrLivArea','BsmtFullBath','BsmtHalfBath','F ullBath', 'HalfBath', 'BedroomAbvGr', \ 'KitchenAbvGr', 'TotRmsAbvGrd', 'Fireplaces', 'GarageYrBlt', 'G arageCars','GarageArea','WoodDeckSF',\ 'OpenPorchSF', 'EnclosedPorch', '3SsnPorch', 'ScreenPorch', 'Po olArea', 'MiscVal', 'MoSold', 'YrSold']

```
In [5]: # preprocess with pipeline and columntransformer
        from sklearn.compose import ColumnTransformer
        from sklearn.pipeline import Pipeline
        from sklearn.preprocessing import OneHotEncoder
        from sklearn.preprocessing import OrdinalEncoder
        from sklearn.preprocessing import StandardScaler
        from sklearn.impute import SimpleImputer
        # one-hot encoder
        categorical_transformer = Pipeline(steps=[
            ('imputer', SimpleImputer(strategy='constant',fill value='missing'
        )),
             ('onehot', OneHotEncoder(sparse=False, handle unknown='ignore'))])
        # ordinal encoder
        ordinal_transformer = Pipeline(steps=[
            ('imputer2', SimpleImputer(strategy='constant',fill_value='NA')),
            ('ordinal', OrdinalEncoder(categories = ordinal cats))])
        # standard scaler
        numeric transformer = Pipeline(steps=[
            ('scaler', StandardScaler())])
        # collect the transformers
        preprocessor = ColumnTransformer(
            transformers=[
                ('num', numeric transformer, num ftrs),
                ('cat', categorical transformer, cat ftrs),
                 ('ord', ordinal transformer, ordinal ftrs)])
In [6]: # fit transform the training set
        X prep = preprocessor.fit transform(X train)
        # little hacky, but collect feature names
        feature names = preprocessor.transformers [0][-1] + \
                        list(preprocessor.named transformers ['cat'][1].get feat
        ure names(cat ftrs)) + \
                        preprocessor.transformers [2][-1]
        df_train = pd.DataFrame(data=X_prep,columns=feature_names)
        print(df train.shape)
        # transform the test
        df test = preprocessor.transform(X_test)
        df test = pd.DataFrame(data=df test,columns = feature names)
        print(df test.shape)
```

```
(876, 221)
(292, 221)
(292, 221)
```

# transform the holdout

print(df\_holdout.shape)

df holdout = preprocessor.transform(X holdout)

df holdout = pd.DataFrame(data=df holdout,columns = feature names)

#### 2) Continuous features: mean or median imputation

- · Imputation means you infer the missing values from the known part of the data
- sklearn's SimpleImputer can do mean and median imputation
- USUALLY A BAD IDEA!
  - MCAR: mean/median of non-missing values is the same as the mean/median of the true underlying distribution, but the variances are different
  - not MCAR: the mean/median and the variance of the completed dataset will be off
  - supervised ML model is too confident (MCAR) or systematically off (not MCAR)

#### 3) Exclude points or features with missing values

- · easy to do with pandas
- it is an ACCEPTABLE approach under two conditions:
  - Little's test supports MCAR (p > 0.05)
  - only small fraction of points contain missing values (maybe a few percent?)
  - the missing values are limited to one or a few features and a large fraction of points are missing from those features (maybe up to 90%?)
- if the MCAR assumption is justified, dropping points will not introduce biases to your model
- · due to the smaller sample size, the confidence of your model might suffer.
- what will you do with missing values when you deploy the model?

```
In [7]: print('data dimensions:',df train.shape)
        print('the p value of the mcar test:', mcar test(df train))
        perc missing per ftr = df train.isnull().sum(axis=0)/df train.shape[0]
        print('fraction of missing values in features:')
        print(perc missing per ftr[perc missing per ftr > 0])
        frac missing = sum(df train.isnull().sum(axis=1)!=0)/df train.shape[0]
        print('fraction of points with missing values:',frac missing)
        data dimensions: (876, 221)
        the p value of the mcar test: 0.029160269814447304
        fraction of missing values in features:
        LotFrontage
                       0.173516
        MasVnrArea
                       0.004566
                       0.050228
        GarageYrBlt
        dtype: float64
        fraction of points with missing values: 0.2237442922374429
```

```
In [8]: print(df_train.shape)
# by default, rows/points are dropped
df_r = df_train.dropna()
print(df_r.shape)
# drop features with missing values
df_c = df_train.dropna(axis=1)
print(df_c.shape)

(876, 221)
(680, 221)
(876, 218)
```

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### XGBoost and missing values

- sklearn raises an error if the feature matrix (X) contains nans.
- · XGBoost doesn't!
- If a feature with missing values is split:
  - XGBoost tries to put the points with missing values to the left and right
  - calculates the impurity measure for both options
  - puts the points with missing values to the side with the lower impurity
- if missingness correlates with the target variable, XGBoost extracts this info!

```
In [9]: import xgboost
        from sklearn.model selection import ParameterGrid
        from sklearn.metrics import mean squared error
        param_grid = {"learning_rate": [0.03],
                      "n estimators": [2000],
                      "seed": [0],
                      #"n jobs": [6],
                      #"reg alpha": [0e0,0.1,0.31622777,1.,3.16227766,10.],
                      #"reg lambda": [0e0,0.1,0.31622777,1.,3.16227766,10.],
                      "missing": [np.nan],
                      #"max depth": [1,2,3,4,5],
                      "colsample_bytree": [0.9],
                      "subsample": [0.66]}
        XGB = xgboost.XGBRegressor()
        XGB.set params(**ParameterGrid(param grid)[0])
        XGB.fit(df train,y train,early stopping rounds=50,eval set=[(df test, y
        test)], verbose=False)
        print('the test RMSE:',XGB.evals result()['validation 0']['rmse'][-1])
        y holdout pred = XGB.predict(df holdout)
        print('the holdout RMSE:',np.sqrt(mean_squared_error(y_holdout,y_holdout
        _pred)))
```

```
/anaconda3/envs/datasci_v0.0.2_local4.yml/lib/python3.6/site-packages/x
gboost/core.py:587: FutureWarning: Series.base is deprecated and will b
e removed in a future version
  if getattr(data, 'base', None) is not None and \
the test RMSE: 24356.306641
the holdout RMSE: 33267.239641714295
```

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### **Multivariate Imputation**

- models each feature with missing values as a function of other features, and uses that estimate for imputation
  - at each step, a feature column is designated as target variable y and the other feature columns are treated as feature matrix X
  - a regressor is trained on (X, y) for known y
  - then, the regressor is used to predict the missing values of y
- in the ML pipeline:
  - create n imputed datasets
  - run all of them through the ML pipeline
  - generate n holdout scores
  - the uncertainty in the holdout scores is due to the uncertainty in imputation
- · works on MCAR and MAR, fails on MNAR
- paper here (https://www.jstatsoft.org/article/view/v045i03)

## sklearn's IterativeImputer

```
LotFrontage MasVnrArea GarageYrBlt
0
     0.424926
                -0.573303
                              0.979398
1
          NaN
                 0.492835
                              1.018748
2
                -0.573303
                              0.192399
          NaN
3
     -0.049970
                 0.810076
                             -0.476551
4
    -1.474659
                -0.022031
                              0.979398
  LotFrontage MasVnrArea GarageYrBlt
0
     0.424926 - 0.573303
                              0.979398
1
    -1.258797
                 0.492835
                              1.018748
    -0.516232
                -0.573303
                              0.192399
3
    -0.049970
                 0.810076
                             -0.476551
    -1.474659
                -0.022031
                              0.979398
```

/anaconda3/envs/datasci\_v0.0.2\_local4.yml/lib/python3.6/site-packages/s klearn/impute/\_iterative.py:599: ConvergenceWarning: [IterativeImputer] Early stopping criterion not reached.

" reached.", ConvergenceWarning)

/anaconda3/envs/datasci\_v0.0.2\_local4.yml/lib/python3.6/site-packages/x gboost/core.py:587: FutureWarning: Series.base is deprecated and will be removed in a future version

if getattr(data, 'base', None) is not None and \

the test RMSE: 24731.107422

the holdout RMSE: 34163.64520140612

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### Reduced-features model (or pattern submodel approach)

- first described in 2007 in a <u>JMLR article (http://www.jmlr.org/papers/v8/saar-tsechansky07a.html)</u> as the reduced features model
- in 2018, "rediscovered" as the pattern submodel approach in <u>Biostatistics</u> (https://www.ncbi.nlm.nih.gov/pubmed/30203058)

#### My holdout set:

index	feature 1	feature 2	feature 3	target var
0	NA	45	NA	0
1	NA	NA	8	1
2	12	6	34	0
3	1	89	NA	0
4	0	NA	47	1
5	687	24	67	1
6	NA	23	NA	1

To predict points 0 and 6, I will use train and test points that are complete in feature 2.

To predict point 1, I will use train and test points that are complete in feature 3.

To predict point 2 and 5, I will use train and test points that are complete in features 1-3.

Etc. We will train as many models as the number of patterns in holdout.

### How to determine the patterns?

```
In [12]: mask = df_holdout[['LotFrontage','MasVnrArea','GarageYrBlt']].isnull()
    unique_rows, counts = np.unique(mask, axis=0,return_counts=True)
    print(unique_rows.shape) # 6 patterns, we will train 6 models
    for i in range(len(counts)):
        print(unique_rows[i],counts[i])
```

```
(6, 3)
[False False False] 223
[False False True] 21
[False True False] 1
[ True False False] 44
[ True False True] 2
[ True True False] 1
```

```
In [13]: def xgb model(X train, Y train, X test, Y test, X holdout, Y holdout, ve
         rbose=1):
             # make into row vectors to avoid an obnoxious sklearn/xgb warning
             Y_train = np.reshape(np.array(Y_train), (1, -1)).ravel()
             Y_test = np.reshape(np.array(Y_test), (1, -1)).ravel()
             Y_holdout = np.reshape(np.array(Y_holdout), (1, -1)).ravel()
             XGB = xgboost.XGBRegressor(n jobs=1)
             # find the best parameter set
             param_grid = {"learning_rate": [0.03],
                            "n estimators": [2000],
                            "seed": [0],
                            #"n jobs": [6],
                            #"reg alpha": [0e0,0.1,0.31622777,1.,3.16227766,10.],
                            #"reg lambda": [0e0,0.1,0.31622777,1.,3.16227766,10.],
                            "missing": [np.nan],
                            #"max depth": [1,2,3,4,5],
                            "colsample bytree": [0.9],
                            "subsample": [0.66]}
             pg = ParameterGrid(param_grid)
             scores = np.zeros(len(pg))
             for i in range(len(pg)):
                 if verbose >= 5:
                      print("Param set " + str(i + 1) + " / " + str(len(pg)))
                 params = pq[i]
                 XGB.set params(**params)
                 eval set = [(X test, Y test)]
                 XGB.fit(X train, Y train,
                         early stopping rounds=50, eval set=eval set, verbose=Fal
         se) # with early stopping
                 Y test pred = XGB.predict(X test, ntree limit=XGB.best ntree lim
         it)
                 scores[i] = mean squared error(Y test,Y test pred)
             best params = np.array(pg)[scores == np.max(scores)]
             if verbose >= 4:
                 print('Test set max score and best parameters are:')
                 print(np.max(scores))
                 print(best params)
             # test the model on the holdout set with best parameter set
             XGB.set params(**best params[0])
             XGB.fit(X train, Y train,
                      early stopping rounds=50, eval set=eval set, verbose=False)
             Y holdout pred = XGB.predict(X holdout, ntree limit=XGB.best ntree l
         imit)
             if verbose >= 1:
                 print ('The MSE is:', mean squared error(Y holdout, Y holdout pred
         ))
             if verbose >= 2:
```

```
print ('The predictions are:')
        print (Y_holdout_pred)
    if verbose >= 3:
        print("Feature importances:")
        print(XGB.feature_importances_)
    return (mean squared error(Y holdout, Y holdout pred), Y holdout pred
, XGB.feature importances )
# Function: Reduced-feature XGB model
# all the inputs need to be pandas DataFrame
def reduced feature xgb(X train, Y train, X test, Y test, X holdout, Y h
oldout):
    # find all unique patterns of missing value in holdout set
    mask = X_holdout.isnull()
    unique rows = np.array(np.unique(mask, axis=0))
    all_Y_holdout_pred = pd.DataFrame()
   print('there are', len(unique rows), 'unique missing value pattern
s.')
    # divide holdout sets into subgroups according to the unique pattern
    for i in range(len(unique rows)):
        print ('working on unique pattern', i)
        ## generate X holdout subset that matches the unique pattern i
        sub X holdout = pd.DataFrame()
        sub Y holdout = pd.Series()
        for j in range(len(mask)): # check each row in mask
            row mask = np.array(mask.iloc[j])
            if np.array_equal(row_mask, unique_rows[i]): # if the patter
n matches the ith unique pattern
                sub X holdout = sub X holdout.append(X holdout.iloc[j])#
append the according X holdout row j to the subset
                sub Y holdout = sub Y holdout.append(Y holdout.iloc[[j
]]) # append the according Y holdout row j
        sub X holdout = sub X holdout[X holdout.columns[~unique rows[i
]]]
        ## choose the according reduced features for subgroups
        sub X train = pd.DataFrame()
        sub Y train = pd.DataFrame()
        sub X test = pd.DataFrame()
        sub Y test = pd.DataFrame()
        # 1.cut the feature columns that have nans in the according sub
X holdout
        sub X train = X train[X train.columns[~unique rows[i]]]
        sub X test = X test[X test.columns[~unique rows[i]]]
        # 2.cut the rows in the sub X train and sub X test that have any
nans
        sub X train = sub X train.dropna()
        sub X test = sub X test.dropna()
        # 3.cut the sub Y train and sub Y test accordingly
        sub Y train = Y train.iloc[sub X train.index]
        sub Y test = Y test.iloc[sub X test.index]
```

```
# run XGB
        sub Y holdout pred = xgb model(sub X train, sub Y train, sub X t
est,
                                       sub Y test, sub X holdout, sub Y
holdout, verbose=0)
        sub_Y_holdout_pred = pd.DataFrame(sub_Y_holdout_pred[1],columns=
['sub Y holdout pred'],
                                          index=sub Y holdout.index)
        print('
                  RMSE:',np.sqrt(mean_squared_error(sub_Y_holdout,sub_Y_
holdout pred)))
        # collect the holdout predictions
        all Y holdout pred = all Y holdout pred.append(sub Y holdout pre
d)
    # rank the final Y holdout pred according to original Y holdout inde
    all Y holdout pred = all Y holdout pred.sort index()
    Y holdout = Y holdout.sort index()
    # get global RMSE
    total RMSE = np.sqrt(mean squared error(Y holdout,all Y holdout pred
))
    return total RMSE
```

#### A python implementation is available on the skipped slide

```
In [14]: print('final RMSE:',reduced_feature_xgb(df_train, y_train, df_test, y_te
         st, df holdout, y holdout))
         there are 6 unique missing value patterns.
         working on unique pattern 0
            RMSE: 36882.33649761256
         working on unique pattern 1
            RMSE: 14122.835714181436
         working on unique pattern 2
            RMSE: 7912.15625
         working on unique pattern 3
            RMSE: 19059.686269128022
         working on unique pattern 4
            RMSE: 20818.556654658543
         working on unique pattern 5
            RMSE: 55023.453125
         final RMSE: 33488.799390845474
```

By the end of this workshop, you will be able to

- Describe the three main types of missingness patterns
- · Evaluate simple approaches for handling missing values
- · Apply XGBoost to a dataset with missing values
- · Apply multivariate imputation
- Apply the reduced-features model (also called the pattern submodel approach)
- · Decide which approach is best for your dataset

### Which approach is best for my data?

- **XGB**: run *n* XGB models with *n* different seeds
- **imputation**: prepare *n* different imputations and run *n* XGB models on them
- **reduced-features**: run *n* reduced-features model with *n* different seeds
- · rank the three methods based on how significantly different the corresponding mean scores are
  - I hope to talk about the results of this experiment at ODSC East next year!

#### Now you can

- Describe the three main types of missingness patterns
- · Evaluate simple approaches for handling missing values
- Apply XGBoost to a dataset with missing values
- Apply multivariate imputation
- Apply the reduced-features model (also called the pattern submodel approach)
- · Decide which approach is best for your dataset

### We are hiring!

- If you don't mind the harsh New England winters and you enjoy working in an academic environment, please come and talk to me or send an email (andras\_zsom@brown.edu), and look out for the job ad at <a href="mailto:Jobs@Brown">Jobs@Brown (https://brown.wd5.myworkdayjobs.com/staff-careers-brown/jobs</a>) (it will be posted in a few days).
- The successful applicant will collaborate with Brown's Advancement, they will work on academic research projects with faculty members, and they will be encouraged to organize and teach at workshops and supervise interns.
- MSc is required!
- PhD and/or industry experience preferred.
- Earliest starting date: February 1st 2020.

### Thanks for your attention!

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TII [ ] •	·		