1 Model Complexity and Model Selection

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1 Question 1 KNN Regressor

import packages

```
In [1]: from scipy.spatial import KDTree
    from sklearn.base import BaseEstimator
    from sklearn.datasets import load_diabetes
    from sklearn.model_selection import train_test_split
    import numpy as np
```

Task I. KNN Regressor Implementation

```
In [2]: class KnnRegressor(BaseEstimator):
    def __init__(self, k=5):
        self.k = k

    def fit(self, X, y):
        self.y_train_ = np.array(y)
        self.x_train_kdtree_ = KDTree(X)
        return self

def predict(self, X):
    # Find k nearest neighbors
    __, neighbours = self.x_train_kdtree_.query(X, k=self.k)
        neighbours = neighbours.reshape(len(X), self.k)

# Get neighbor labels (regression targets here)
        neighbour_values = self.y_train_[neighbours]

# Compute mean
    y_pred = neighbour_values.mean(axis=1)
```

Task II. Test implementation

Load dataset

```
In [3]: # Load dataset
    diabetes = load_diabetes()
    X = diabetes.data
    y = diabetes.target
    X.shape, y.shape
Out[3]: ((442, 10), (442,))
```

Split dataset

Guess a K

According to the empirical rule, $\sqrt{trainingSetSize}$ is often close to the optimal value of K. Therefore, I estimate that K = $\sqrt{309} \approx$ 18 is an acceptable choice for K.

Test

```
In [5]: # Test KNN Regressor
knn = KnnRegressor(k=18)
knn.fit(X_train, y_train)
y_hat_train = knn.predict(X_train)
y_hat_test = knn.predict(X_test)
y_hat_train.shape, y_hat_test.shape
```

```
Out[5]: ((309,), (133,))
```

Evaulation

Use the sum of the squares of the errors as the error function to measure training errors and testing errors.

```
In [6]: # calculate the sum of squared errors
def sse(y_true, y_pred):
    return np.sum((y_true - y_pred) ** 2)
```

Compute error

When K = 18 The training error is 873984.8395061729 The testing error is 468777.64197530865

Further test to find optimal K

Test and find the optimal K (1~30)

```
sse_test.append(sse(y_test, y_hat_test))
return sse_train, sse_test

sse_train, sse_test = choose_k(X_train, y_train, X_test, y_test, 30)

# find the minimum sse_test and the corresponding k
min_sse_test = min(sse_test)
best_k = sse_test.index(min_sse_test) + 1
print(f"The minimum testing error is {min_sse_test} when K = {best_k}")
print(f"Hence, the optimal K is {best_k}")
```

The minimum testing error is 450863.84171597636 when K = 26 Hence, the optimal K is 26

2 Question 2 L-fold Cross Validation

Task I. L-fold Implementation

```
In [9]: class LFold:
            def init (self, L=5, random state=None, shuffle=False):
                 self.L = L
                 self.random state = random state
                 self.shuffle = shuffle
            def get_n_splits(self):
                 return self.L
            def split(self, x, y=None, groups=None):
                 n \text{ samples} = len(x)
                indices = np.arange(n samples) # indexes of samples
                 if self.shuffle:
                     rng = np.random.RandomState(self.random_state)
                     rnq.shuffle(indices)
                # calculate the size of each fold
                fold_size = np.full(self.L, n_samples // self.L, dtype=int)
                fold_size[: n_samples % self.L] += 1
                 current = 0
```

```
for f_size in fold_size:
    start, end = current + f_size
    test_idx = indices[start:end]
    train_idx = np.concatenate([indices[:start], indices[end:]])
    yield train_idx, test_idx
    current = end
```

Test the implementation

```
In [10]: for idx_train, idx_test in LFold(L = 5).split(list(range(20))):
    print(idx_train, idx_test)

[ 4  5  6  7  8  9 10 11 12 13 14 15 16 17 18 19] [0 1 2 3]
    [ 0  1  2  3  8  9 10 11 12 13 14 15 16 17 18 19] [4  5  6  7]
    [ 0  1  2  3  4  5  6  7 12 13 14 15 16 17 18 19] [8  9 10 11]
    [ 0  1  2  3  4  5  6  7  8  9 10 11 16 17 18 19] [12 13 14 15]
    [ 0  1  2  3  4  5  6  7  8  9 10 11 12 13 14 15] [16 17 18 19]
```

Task II. Test the effect of the k

Test k values from 1 to 60, calculate and compare the mean and standard deviation of MSE for both training and test sets, and evaluate which k value is optimal.

Construct test

Among the 442 samples, a 7:3 split results in 309 training samples. Given the moderate sample size, we use 10-fold cross-validation, so that each training uses about 278 samples and validates on 31 samples. Compared to 5-fold, 10-fold cross-validation provides a more stable estimate of generalization error while keeping the computational cost acceptable.

```
In [11]: import pandas as pd

def mse(y_true, y_pred):
    """Calculate the mean squared error."""
    return np.mean((y_true - y_pred) ** 2)

def cross_validate_knn_lfold(
    X,
    y,
    max_k,
```

```
L=10.
    random state=18.
    shuffle=False,
):
    # initalize the result lists
    train mse means = []
    train mse stds = []
    valid mse means = []
    valid mse stds = []
    # cross validation for each k
    for k in range(1, max k + 1):
       train mses = []
        valid_mses = []
        lfold = LFold(L, shuffle=shuffle, random_state=random_state)
        fold splits = lfold.split(X)
        # iterate over each fold
        for train_idx, valid_idx in fold_splits:
            # split data
            X_tr, y_tr = X[train_idx], y[train_idx]
            X_val, y_val = X[valid_idx], y[valid_idx]
            # train KNN model
            knn = KnnRegressor(k=k)
            knn.fit(X_tr, y_tr)
            # calculate the training set MSE
            y_tr_pred = knn.predict(X_tr)
            train_mse = mse(y_tr, y_tr_pred)
            train mses.append(train mse)
            # calculate the validation set MSE
            v val pred = knn.predict(X val)
            valid_mse = mse(y_val, y_val_pred)
            valid mses.append(valid mse)
        # calculate the mean and standard deviation
       train mse mean = np.mean(train mses)
        train mse std = np.std(train mses)
        valid mse mean = np.mean(valid mses)
        valid_mse_std = np.std(valid_mses)
        train_mse_means.append(train_mse_mean)
```

Run test

Execute cross_validate_knn_lfold function

Out [12]: k_values train_mse_means train_mse_stds valid_mse_means valid_mse_stds

0	1	0.000000	0.000000	5446.774745	851.363011
1	2	1529.635023	66.820172	4225.125843	496.874615
2	3	2022.684618	63.614753	4033.731069	446.768688
3	4	2257.827505	90.719343	3667.087510	211.074126
4	5	2439.041247	25.943655	3620.916153	286.613920

Find optimal k

The k with the smallest valid_mse_means is the optimal k

The optimal K is: 19, the corresponding validation set MSE mean is: 3172.3750

Task III. Analyse the effect of k and L

In this section, I will first test L = 5 and 10 for $k(1\sim60)$ and store the results.

Then, compute the confidence intervals for each result using the following rule:

$$m\pm 1.96s\sqrt{L}$$

Then, I will draw two plots for comparison.

Generate results for L=5 and 10

```
In [14]: L1 = 5
    L2 = 10
    max_k = 60

results_l5 = cross_validate_knn_lfold(
    X,
    y,
    max_k=max_k,
    L=L,
    random_state=18,
    shuffle=True,
)
results_l10 = cross_validate_knn_lfold(
    X,
    y,
    max_k=max_k,
    L=L,
```

```
random_state=18,
    shuffle=True,
)
```

Compute confidence intervals

```
In [15]: means_l5 = results_l5["valid_mse_means"]
    stds_l5 = results_l10["valid_mse_stds"]
    means_l10 = results_l10["valid_mse_means"]
    stds_l10 = results_l10["valid_mse_stds"]

    ci_lowers_l5 = means_l5 - 1.96 * stds_l5 / np.sqrt(L1)
        ci_uppers_l5 = means_l5 + 1.96 * stds_l5 / np.sqrt(L1)
        ci_lowers_l10 = means_l10 - 1.96 * stds_l10 / np.sqrt(L2)
        ci_uppers_l10 = means_l10 + 1.96 * stds_l10 / np.sqrt(L2)

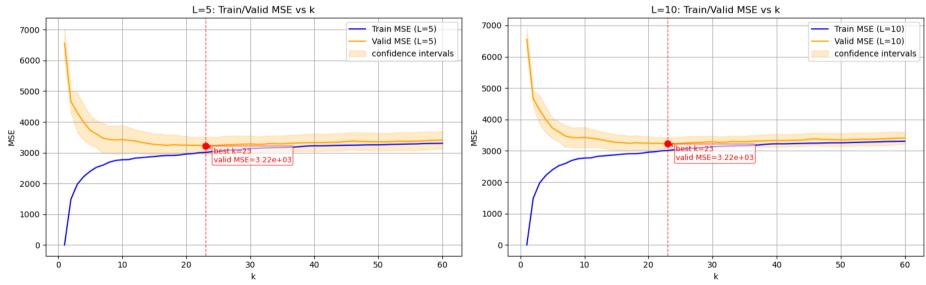
# add to results in order to plot
    results_l5["ci_lowers"] = ci_lowers_l5
    results_l5["ci_uppers"] = ci_uppers_l5
    results_l10["ci_uppers"] = ci_uppers_l10
    results_l10["ci_uppers"] = ci_uppers_l10
```

Draw plots

Plot two figures: one for L=5 and one for L=10. The confidence intervals are represented by shaded areas around the validation MSE mean curves.

```
results 15["ci lowers"],
    results 15["ci uppers"],
    color="orange",
    alpha=0.2.
    label="confidence intervals"
axes[0].scatter(best k l5, best mse l5, color="red", s=60, zorder=5)
axes[0].axvline(best k l5, color="red", linestyle="--", linewidth=1, alpha=0.7)
axes[0].annotate(
   f"best k={best k l5}\nvalid MSE={best mse l5:.3g}",
    xy=(best_k_l5, best_mse_l5),
    xytext=(10, -20),
    textcoords="offset points",
    color="red",
    fontsize=9.
    bbox=dict(boxstyle="round.pad=0.2", fc="white", ec="red", alpha=0.6),
    arrowprops=dict(arrowstyle="->", color="red", lw=1),
axes[0].set xlabel("k")
axes[0].set ylabel("MSE")
axes[0].set title("L=5: Train/Valid MSE vs k")
axes[0].legend()
axes[0].grid(True)
# L=10
best idx l10 = int(np.argmin(results l10["valid mse means"]))
best k l10 = int(k values[best idx l10])
best_mse_l10 = float(results_l10["valid_mse_means"][best idx l10])
axes[1].plot(k values, results l10["train mse means"], label="Train MSE (L=10)", color="blue")
axes[1].plot(k values, results l10["valid mse means"], label="Valid MSE (L=10)", color="orange")
axes[1].fill between(
    k values,
    results l10["ci lowers"],
    results l10["ci uppers"],
    color="orange",
    alpha=0.2,
    label="confidence intervals"
axes[1].scatter(best k l10, best mse l10, color="red", s=60, zorder=5)
axes[1].axvline(best k l10, color="red", linestyle="--", linewidth=1, alpha=0.7)
```

```
axes[1].annotate(
   f"best k={best_k_l10}\nvalid MSE={best_mse_l10:.3g}",
    xy=(best_k_l10, best_mse_l10),
   xytext=(10, -20),
    textcoords="offset points",
    color="red",
    fontsize=9.
    bbox=dict(boxstyle="round,pad=0.2", fc="white", ec="red", alpha=0.6),
    arrowprops=dict(arrowstyle="->", color="red", lw=1),
axes[1].set xlabel("k")
axes[1].set ylabel("MSE")
axes[1].set_title("L=10: Train/Valid MSE vs k")
axes[1].legend()
axes[1].grid(True)
plt.tight layout()
plt.show()
```



Analysis

From the plots, we observe that very small values of k (e.g., 1–3) lead to extremely low training errors but high validation errors, indicating overfitting. When k increases to 23, the validation error reaches its minimum, which can be regarded as the optimal region. For large k

values (>23), both training and validation errors increase, and the model suffers from underfitting.

Comparing L=5 and L=10, both curves show similar trends, but the confidence intervals are wider for L=5, implying higher variance in the error estimates. In contrast, L=10 produces narrower confidence intervals and thus more stable performance estimates, at the cost of higher computational effort. Given the dataset size (309 training samples), L=10 provides a more reliable choice.

3 Question 3 Automatic Model Selection

Task I. KnnRegressorCV implementation

```
In [17]: from sklearn.base import BaseEstimator
         class KnnRegressorCV(BaseEstimator):
             def __init__(self, ks=list(range(1, 21)), cv=LFold(5)):
                  self.ks = ks
                 self.cv = cv
                 self.k_ = None # the best k
                  self.model = None # the best model
             def fit(self, x, y):
                  best k = None
                 best_score = float("inf")
                 # iterate over each k
                 for k in self.ks:
                     val errors = []
                     for train_idx, val_idx in self.cv.split(x):
                         # split data
                         X_tr, y_tr = x[train_idx], y[train_idx]
                         X_{val}, y_{val} = x[val_idx], y[val_idx]
                          knn = KnnRegressor(k=k)
                         knn.fit(X_tr, y_tr)
                         y_val_pred = knn.predict(X_val)
                         val_mse = mse(y_val, y_val_pred)
                         val_errors.append(val_mse)
                     mean_error = np.mean(val_errors)
```

```
if mean_error < best_score:
    best_score = mean_error
    best_k = k

# record the best k and the model
self.k_ = int(best_k)
self.model_ = KnnRegressor(k=self.k_).fit(x, y)
return self

def predict(self, x):
    return self.model_.predict(x)</pre>
```

```
In [18]: l_fold = LFold(L=10)
knn_cv = KnnRegressorCV(ks=list(range(1, 21)), cv=l_fold)
knn_cv.fit(X, y)
knn_cv.predict(X_test)
print("Chosen k by inner CV:", knn_cv.k_)
Chosen k by inner CV: 19
```

Task II. Nested cross-validation

Experimental setup: Outer CV is 10-fold, inner CV is 10-fold, and the tested k values are in the range [1, 60]

```
In [19]: outer_cv = LFold(L=10, shuffle=True, random_state=18)
    inner_cv = LFold(L=10, shuffle=True, random_state=18)

    chosen_ks, test_mses = [], []
    fold_count = 0
    for tr_idx, val_idx in outer_cv.split(X):
        X_tr, y_tr = X[tr_idx], y[tr_idx]
        X_val, y_val = X[val_idx], y[val_idx]

        knn = KnnRegressorCV(ks=list(range(1,61)), cv=inner_cv).fit(X_tr, y_tr)
        chosen_ks.append(knn.k_)
        y_pred = knn.predict(X_val)
        # error of best model (in this outer fold)
        test_mses.append(mse(y_val, y_pred))
        fold_count += 1
        print(f"[Outer Fold] {fold_count} completed, Best k: {knn.k_} with Test MSE: {test_mses[-1]:.4f}")

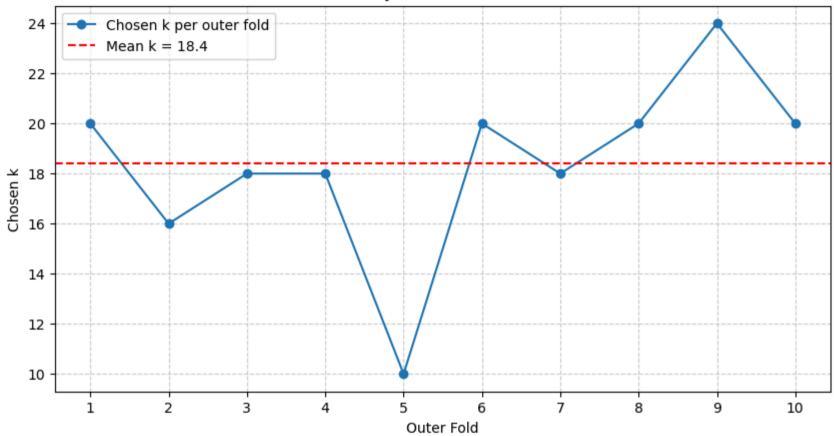
# mean K
```

```
mean k = np.mean(chosen ks)
print(f"Average k: {mean k:.2f}")
# plot the chosen k
plt.figure(figsize=(10, 5))
plt.plot(range(1, len(chosen ks) + 1), chosen ks, marker='o', linestyle='-', label="Chosen k per outer fold")
# mean k
plt.axhline(y=mean k, color='r', linestyle='--', label=f"Mean k = {mean k:.1f}")
plt.xlabel("Outer Fold")
plt.ylabel("Chosen k")
plt.title("K values chosen by Internal CV across Outer Folds")
plt.xticks(range(1, len(chosen_ks) + 1))
plt.legend()
plt.grid(True, linestyle="--", alpha=0.6)
plt.show()
[Outer Fold] 1 completed, Best k: 20 with Test MSE: 3702.6724
[Outer Fold] 2 completed, Best k: 16 with Test MSE: 3814.6490
[Outer Fold] 3 completed, Best k: 18 with Test MSE: 3032.5382
[Outer Fold] 4 completed, Best k: 18 with Test MSE: 3377.8617
[Outer Fold] 5 completed, Best k: 10 with Test MSE: 3465.1839
[Outer Fold] 6 completed, Best k: 20 with Test MSE: 3184.8859
[Outer Fold] 7 completed, Best k: 18 with Test MSE: 2658.6661
[Outer Fold] 8 completed, Best k: 20 with Test MSE: 4207.0656
[Outer Fold] 9 completed, Best k: 24 with Test MSE: 2329.0017
```

[Outer Fold] 10 completed, Best k: 20 with Test MSE: 2886.2298

Average k: 18.40

K values chosen by Internal CV across Outer Folds



Analysis

The KNN regressor with internal cv selected an average k-value of 18.4 across the 10 outer folds. The individual k-values chosen ranged from 10 to 24, with most selections clustering around 18-20.

However, the mean k-value of 18.4 does **NOT** correspond to the optimal k-value with respect to the outer test sets. The lowest test MSE (2329.0017) was achieved in outer fold 9 with k=24.

First, the representativeness and stability of the dataset will affect the selection of the optimal model. We can see that in this experiment, the k values chosen by different outer folds vary greatly (10-24), which shows that there are large differences within the dataset. When the training subsets cannot well represent the overall data distribution, the selection by internal CV becomes unreliable.

Second, the choice of hyperparameters will also affect model selection. KNN is very sensitive to the k value, and small changes may lead to large performance fluctuations. If the model has many "local optimal points", internal CV can easily get trapped in them.

In addition, the design of cross-validation, such as the number of folds, splitting method, and random seed, will also affect the stability of the results.

In []: