a) The scientist believes that past rainfall and streamflow data (with lags up to 5 days) may be important, but they are unsure which specific lag times are most relevant for each variable. Additionally, they want to avoid using all lagged variables to keep the number of model parameters manageable. If they want to build a model with 5 inputs, how can they determine which lagged variables to include in the model? [Provide a written explanation—no coding or calculations are needed].

In my opinion, I think they could calculate the correlations between the streamflow and the lagged variables. This way will help them to identify which lags have the strongest relationship with streamflow. Also, they could test different combinations of lagged variables through calculate the AIC. And this scientist are familiar with knowledge about hydrological processes, they could determine which lagged variables to include in the model based on their experience.

b) Build a multiple linear regression model to predict streamflow using the following five input variables: rainfall at lag time ($\tau = 1, 5$) and streamflow at lag times ($\tau = 1, 3, 5$):

```
In [97]: import pandas as pd
import numpy as np
from sklearn.linear_model import LinearRegression
from sklearn.metrics import mean_squared_error, r2_score
from scipy.stats import pearsonr
import matplotlib.pyplot as plt
```

```
In [98]: # Load dataset
df = pd.read_csv('fish_river.csv')
df
```

Out[98]:		datetime	year	rainfall	runoff
	0	2016-01-01	2016	0.000675	1720.0
	1	2016-01-02	2016	0.144395	1680.0
	2	2016-01-03	2016	2.918217	1640.0
	3	2016-01-04	2016	0.000000	1580.0
	4	2016-01-05	2016	0.000000	1520.0
	•••	•••			
	2187	2021-12-27	2021	0.000000	989.0
	2188	2021-12-28	2021	3.711825	972.0
	2189	2021-12-29	2021	0.018728	947.0
	2190	2021-12-30	2021	0.000000	928.0
	2191	2021-12-31	2021	0.030013	902.0

2192 rows × 4 columns

```
In [99]: # Create lag variables for rainfall and streamflow
for lag in [1, 3, 5]:
    df[f'runoff_lag_{lag}'] = df['runoff'].shift(lag)
for lag in [1, 5]:
    df[f'rainfall_lag_{lag}'] = df['rainfall'].shift(lag)
df = df.dropna()
df
```

Out[99]:		datetime	year	rainfall	runoff	runoff_lag_1	runoff_lag_3	runoff_lag_5	ra
	5	2016-01- 06	2016	0.000000	1450.0	1520.0	1640.0	1720.0	
	6	2016-01- 07	2016	0.000000	1390.0	1450.0	1580.0	1680.0	
	7	2016-01- 08	2016	0.000000	1350.0	1390.0	1520.0	1640.0	
	8	2016-01- 09	2016	1.009949	1310.0	1350.0	1450.0	1580.0	
	9	2016-01- 10	2016	19.440051	1290.0	1310.0	1390.0	1520.0	
	•••						•••		
	2187	2021-12- 27	2021	0.000000	989.0	1020.0	1120.0	1120.0	
	2188	2021-12- 28	2021	3.711825	972.0	989.0	1070.0	1130.0	
	2189	2021-12- 29	2021	0.018728	947.0	972.0	1020.0	1120.0	
	2190	2021-12- 30	2021	0.000000	928.0	947.0	989.0	1070.0	
	2191	2021-12- 31	2021	0.030013	902.0	928.0	972.0	1020.0	

2187 rows × 9 columns

```
In [100... # get the training data and testing data
    train_data = df[(df['year'] >= 2016) & (df['year'] <= 2019)]
    test_data = df[(df['year'] >= 2020) & (df['year'] <= 2021)]

# define data
X_train = train_data[['runoff_lag_1', 'runoff_lag_3', 'runoff_lag_5', 'rainf
y_train = train_data['runoff']
X_test = test_data[['runoff_lag_1', 'runoff_lag_3', 'runoff_lag_5', 'rainfal
y_test = test_data['runoff']
X_train</pre>
```

Out[100...

	runoff_lag_1	runoff_lag_3	runoff_lag_5	rainfall_lag_1	rainfall_lag_5
5	1520.0	1640.0	1720.0	0.000000	0.000675
6	1450.0	1580.0	1680.0	0.000000	0.144395
7	1390.0	1520.0	1640.0	0.000000	2.918217
8	1350.0	1450.0	1580.0	0.000000	0.000000
9	1310.0	1390.0	1520.0	1.009949	0.000000
•••	•••	•••	•••		
1456	1630.0	1760.0	1900.0	0.000000	3.488331
1457	1570.0	1690.0	1830.0	4.640274	0.006127
1458	1510.0	1630.0	1760.0	0.000000	0.000000
1459	1450.0	1570.0	1690.0	0.000000	0.000000
1460	1400.0	1510.0	1630.0	2.977127	0.000000

1456 rows × 5 columns

In [101... X_test

Out[101...

	runoff_lag_1	runoff_lag_3	runoff_lag_5	rainfall_lag_1	rainfall_lag_5
1461	1350.0	1450.0	1570.0	7.502306	4.640274
1462	1300.0	1400.0	1510.0	0.440019	0.000000
1463	1250.0	1350.0	1450.0	0.060312	0.000000
1464	1200.0	1300.0	1400.0	0.000000	2.977127
1465	1160.0	1250.0	1350.0	0.668274	7.502306
•••	•••	•••		•••	•••
2187	1020.0	1120.0	1120.0	0.000000	14.028154
2188	989.0	1070.0	1130.0	0.000000	0.000000
2189	972.0	1020.0	1120.0	3.711825	0.000000
2190	947.0	989.0	1070.0	0.018728	0.000000
2191	928.0	972.0	1020.0	0.000000	0.000000

731 rows × 5 columns

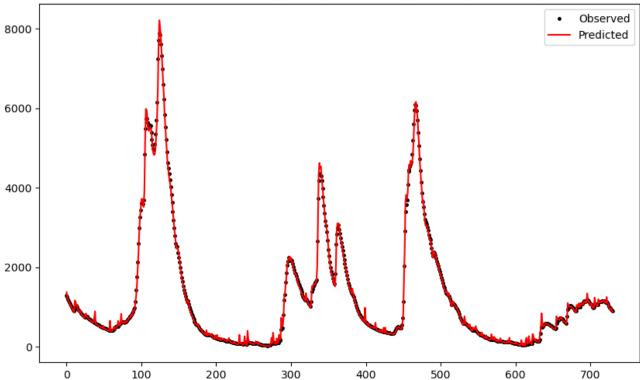
In [102... # Initialize the multiple linear regression model model = LinearRegression() model.fit(X train, y train) # Make predictions y_pred = model.predict(X_test) # compute the RMSE rmse = np.sqrt(mean_squared_error(y_test, y_pred)) print(f'RMSE: {rmse}') # compute the Pearson correlation coefficient pearson_corr, _ = pearsonr(y_test, y_pred) print(f'Pearson Correlation Coefficient: {pearson_corr}') # compute the relative bias relative_bias = (y_pred.mean() - y_test.mean()) / y_test.mean() print(f'Relative Bias: {relative_bias}') # Plot the model predictions for the testing period (2020-2021) plt.figure(figsize=(10, 6)) plt.plot(obs, 'k.', label='Observed', markersize=5) plt.plot(y_pred, 'r-', label='Predicted') plt.title('Runoff Prediction using Multiple Linear Regression (2020-2021)') plt.legend() plt.show()

RMSE: 98.10814034558912

Pearson Correlation Coefficient: 0.9979102416958303

Relative Bias: -0.00023792941723045185

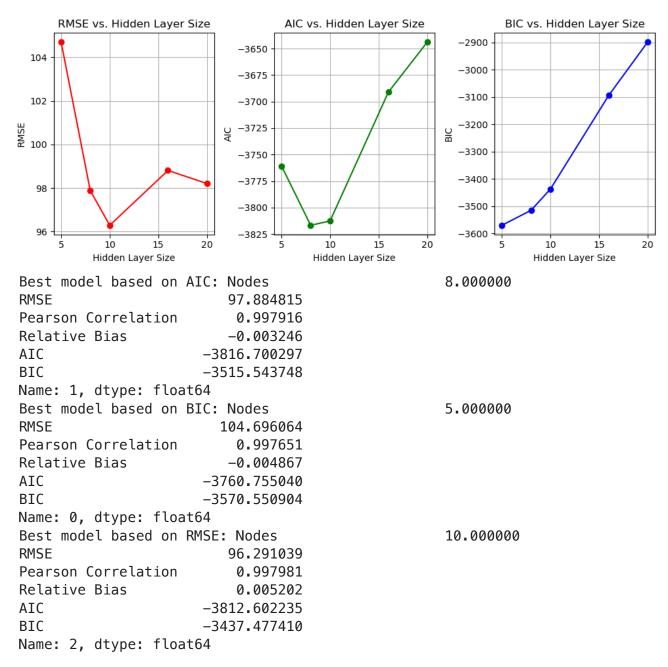
Runoff Prediction using Multiple Linear Regression (2020-2021)



c) Build a Multilayer Perceptron (MLP) model to predict streamflow using the same five input variables in (b) and the same training-testing data split. [Hint: For Python users, utilize the Scikit-Learn MLPRegressor model with default settings unless otherwise specified]

```
In [103... | nodes_list = [5, 8, 10, 16, 20]
         results = []
         for nodes in nodes list:
             # # Define the MLP model
             mlp = MLPRegressor(hidden_layer_sizes=(nodes,), max_iter=10000, random_s
             # fit the model
             mlp.fit(X_train, y_train)
             # get the y_predict
             y_pred = mlp.predict(X_test)
             # compute the RMSE
             rmse = np.sqrt(mean_squared_error(y_test, y_pred))
             # compute Pearson
             pearson_corr, _ = pearsonr(y_test, y_pred)
             # compute the relative bias
             relative_bias = (y_pred.mean() - y_test.mean()) / y_test.mean()
             # compute the number of parameters
             K = (nodes * 5) + (1 * nodes) + nodes + 1
             # compute the AIC and BIC
             n = len(y_train)
             aic = 2 * K + n * np.log(rmse/n)
             bic = K * np.log(n) + n * np.log(rmse/n)
             results.append((nodes, rmse, pearson_corr, relative_bias, aic, bic))
         # transform the results to DataFrame
          results_df = pd.DataFrame(results, columns=['Nodes', 'RMSE', 'Pearson Correl
         print(results_df)
         # plot AIC, BIC and RMSE
         fig, axs = plt.subplots(1, 3, figsize=(10, 4))
         axs[0].plot(results df['Nodes'], results df['RMSE'], marker='o', color='red'
         axs[0].set_title('RMSE vs. Hidden Layer Size')
         axs[0].set_xlabel('Hidden Layer Size')
```

```
axs[0].set_ylabel('RMSE')
 axs[0].grid()
 axs[1].plot(results_df['Nodes'], results_df['AIC'], marker='o', color='greer
 axs[1].set_title('AIC vs. Hidden Layer Size')
 axs[1].set xlabel('Hidden Layer Size')
 axs[1].set ylabel('AIC')
 axs[1].grid()
 axs[2].plot(results_df['Nodes'], results_df['BIC'], marker='o', color='blue'
 axs[2].set_title('BIC vs. Hidden Layer Size')
 axs[2].set_xlabel('Hidden Layer Size')
 axs[2].set ylabel('BIC')
 axs[2].grid()
 plt.tight layout()
 plt.show()
 # find the best model
 best_aic_model = results_df.loc[results_df['AIC'].idxmin()]
 best_bic_model = results_df.loc[results_df['BIC'].idxmin()]
 best rmse model = results df.loc[results df['RMSE'].idxmin()]
 print(f'Best model based on AIC: {best aic model}')
 print(f'Best model based on BIC: {best_bic_model}')
 print(f'Best model based on RMSE: {best_rmse_model}')
   Nodes
                RMSE Pearson Correlation Relative Bias
                                                                   AIC
0
       5
         104.696064
                                 0.997651
                                                -0.004867 -3760.755040
1
       8
           97.884815
                                 0.997916
                                                -0.003246 -3816.700297
2
      10
           96.291039
                                 0.997981
                                                 0.005202 -3812.602235
3
                                                -0.003709 -3691.058442
      16
           98.806248
                                 0.997901
4
      20
           98.208742
                                 0.997977
                                                 0.012131 -3643.889957
           BIC
0 -3570.550904
1 -3515.543748
2 -3437.477410
3 -3094.028792
4 -2898.923757
```



Do any of the five MLP models outperform the multiple-linear regression model? If so, why might that be the case?

Yes. MLP can capture nonlinear relationships between the input and output variables better than MLR. So if the relationship is nonlinear, MLP may provide better predictions. And MLP especially with more nodes and layers, have a higher capacity to learn complex patterns from the data. However, if an MLP model is too complex relative to the size of the training dataset, it may overfit the training data. Also, if the dataset is small, linear regression might perform better.

Identify the best MLP model based on each metric. Do all metrics agree? If not, explain

why?

No, not all metrics agree. This is because the RMSE foucus on minimizing error, while AIC and BIC foucus on model complexity. So based on AIC and BIC, we could choose the simpler model which RMSE helps us to choose more accuracy model.

The Maine Department of Environmental Protection is particularly interested in developing a data-driven model that excels at predicting extreme runoff values (e.g., the maximum daily runoff in a year or daily runoff values above the 95th percentile). Discuss two approaches that could be used to enhance the model's performance in predicting extreme events, even if this comes at the expense of the model's accuracy in predicting non-extreme runoff values. [Hint: Consider the choice of loss function and regularization techniques.]

1. Loss Functions

Traditional loss functions such as the mean square error (MSE) tend to focus on minimizing the overall error of all predictions, which can lead to poor performance when predicting extreme values. In order to improve the prediction of extreme runoff values, greater weights can be assigned to the extreme values. For example, a weighted loss function could assign a greater penalty to errors in runoff values above the 95th percentile, ensuring that the model prioritizes accurate prediction of high runoff events. Another approach is to use quantile regression, which predicts a specific percentile (e.g., the maximum daily runoff in a year or daily runoff values above the 95th percentile). The quantile loss function minimizes the error for a given quantile, ensuring that the model is appropriate for predicting extreme high runoff events rather than focusing on the median or mean. This helps improve the performance of extreme runoff prediction.

2. regularization techniques

L1 regularization (Lasso) expresses the sparsity of the model, which means that only the most important features for predicting extreme values are retained. Regularization based on Extreme Value Theory (EVT) is specifically designed to improve the prediction of rare and extreme events. By introducing EVT, models can emphasize the tail behavior of the runoff distribution and improve the model's ability to accurately predict large runoff values.