20250425 01

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[1]: # What is Cross-Validation?
      import pandas as pd
      from sklearn.datasets import load_breast_cancer
      from sklearn.model selection import cross val score
      from sklearn.linear_model import LogisticRegression
      from sklearn.tree import DecisionTreeClassifier
      # Loading dataset
      data = load_breast_cancer()
      X = pd.DataFrame(data.data, columns = data.feature_names)
      y = pd.Series(data.target)
[13]: # Creating models
      logreg = LogisticRegression(solver = 'liblinear')
      tree = DecisionTreeClassifier(random_state = 42)
[15]: # Comparing models
      logreg_score = cross_val_score(logreg, X, y, cv = 5)
      print('Logistic Regression cross-validation scores:', logreg_score)
      print('Mean accuracy:', logreg_score.mean())
      tree_score = cross_val_score(tree, X, y, cv = 5)
      print('Decision Tree cross-validation scores:', tree_score)
      print('Mean accuracy:', tree_score.mean())
     Logistic Regression cross-validation scores: [0.92982456 0.93859649 0.97368421
     0.94736842 0.96460177]
     Mean accuracy: 0.9508150908244062
     Decision Tree cross-validation scores: [0.9122807 0.90350877 0.92982456
     0.95614035 0.88495575]
     Mean accuracy: 0.9173420276354604
[23]: # Making pipelines
      from sklearn.pipeline import Pipeline
      from sklearn.preprocessing import StandardScaler
      # Standizing for Logistic Regression, but no need for Decision Tree
```

Logistic Regression Pipeline cross-validation Scores: [0.98245614 0.97368421 0.97368421 0.97368421 0.99115044]

Mean Accuracy: 0.9789318428815402

Decision Tree Pipeline cross-validation Scores: [0.9122807 0.90350877 0.92982456 0.95614035 0.88495575]

Mean Accuracy: 0.9173420276354604

0.0.1 Conclusion:

- Logistic Regression achieved higher and more stable accuracy (mean 0.951) compared to Decision Tree (mean 0.917).
- After introducing a Pipeline with StandardScaler, Logistic Regression performance further improved, confirming its sensitivity to feature scaling.
- Decision Tree was also wrapped in a Pipeline for consistency, though scaling is not needed for tree-based models.