

Model Optimization and Tuning Phase Template

Date	20 July 2024
Team ID	xxxxxxx
Project Title	Detection of Autistic Spectrum Disorder: Classification
Maximum Marks	10 Marks

Model Optimization and Tuning Phase

The Model Optimization and Tuning Phase involves refining neural network models for peak performance. It includes optimized model code, fine-tuning hyperparameters, comparing performance metrics, and justifying the final model selection for enhanced predictive accuracy and efficiency.

Hyperparameter Tuning Documentation (8 Marks):

Model	Tuned Hyperparameters
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Logistic Regression	<p>C: Regularization strength Solver: Optimization algorithm</p> <p>C: Controls the regularization strength. Smaller values specify stronger regularization. Solver: Algorithm to use for optimization.</p> <pre>from sklearn.linear_model import LogisticRegression lgr=LogisticRegression() lgr.fit(X_train,y_train) * LogisticRegression LogisticRegression() pred_lgr=predict(X_test) y_pred_lgr = lgr.predict(X_test) from sklearn.metrics import classification_report accuracy_lgr = accuracy_score(y_test,y_pred_lgr) print('Accuracy LGR:', accuracy_lgr*100) accuracy_lgr = accuracy_score(y_test,y_pred_lgr) print("Accuracy LGR:", accuracy_lgr*100) Accuracy LGR: 100.0 print(classification_report(y_true=y_test,y_pred=pred))</pre> <table><thead><tr><th></th><th>precision</th><th>recall</th><th>f1-score</th><th>support</th></tr></thead><tbody><tr><td>0</td><td>1.00</td><td>1.00</td><td>1.00</td><td>132</td></tr><tr><td>1</td><td>1.00</td><td>1.00</td><td>1.00</td><td>51</td></tr><tr><td>accuracy</td><td></td><td></td><td>1.00</td><td>183</td></tr><tr><td>macro avg</td><td>1.00</td><td>1.00</td><td>1.00</td><td>183</td></tr><tr><td>weighted avg</td><td>1.00</td><td>1.00</td><td>1.00</td><td>183</td></tr></tbody></table>		precision	recall	f1-score	support	0	1.00	1.00	1.00	132	1	1.00	1.00	1.00	51	accuracy			1.00	183	macro avg	1.00	1.00	1.00	183	weighted avg	1.00	1.00	1.00	183
	precision	recall	f1-score	support																											
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weighted avg	1.00	1.00	1.00	183																											
SVM	<p>C: Regularization parameter Kernel: Type of kernel used Gamma: Kernel coefficient</p>																														

	<h2>SVC</h2> <pre>from sklearn.svm import SVC svm=SVC(kernel='rbf', random_state=0) svm.fit(X_train, y_train)</pre> <div> <div>SVC</div> <div>SVC(random_state=0)</div> </div> <pre>y_pred_svc=svm.predict(X_test)</pre> <div>+ Code + Markdown</div> <pre>print('Training Set: ', svm.score(X_train,y_train)) print('Testing Set:',svm.score(X_test,y_test))</pre> <p>Training Set: 0.9530516431924883 Testing Set: 0.9453551912568307</p> <pre>accuracy_SVC=svm.score(X_test,y_test) print('Accuracy_SVM:', accuracy_SVC*100)</pre>
Decision Tree	<p>Max Depth: Maximum depth of the tree Min Samples Split: Minimum number of samples required to split an internal node</p>

	<h2>Decision Tree</h2> <pre> dt = DecisionTreeClassifier() dt.fit(X_train,y_train)] ▼ DecisionTreeClassifier DecisionTreeClassifier()] y_pred_dt=dt.predict(X_test)] print('Training Set: ',dt.score(X_train,y_train)) print('Test Set: ',dt.score(X_test,y_test))] Training Set: 1.0 Test Set: 1.0] print("Accuracy:", metrics.accuracy_score(y_test, y_pred_dt)*100)] Accuracy: 100.0] accuracy_dt=accuracy_score(y_test,y_pred_dt) print('Accuracy DT:', accuracy_dt*100)] </pre>
Random Forest	<p>n_estimators: Number of trees Max Features: Number of features to consider for splitting a node</p>

	<h2>Random Forest</h2> <pre> rand_forest = RandomForestClassifier(random_state=42) rand_forest.fit(X_train, y_train) RandomForestClassifier RandomForestClassifier(random_state=42) predictionRF = rand_forest.predict(X_test) print("Training set: ",rand_forest.score(X_train, y_train)) print("Testing set: ",rand_forest.score(X_test, y_test)) Training set: 1.0 Testing set: 1.0 accuracy_RF=rand_forest.score(X_test, y_test) print ("Accuracy_RF:",accuracy_RF*100) Accuracy_RF: 100.0 </pre>
KNN	<p>n_neighbors: Number of neighbors Metric: Distance metric</p> <h2>KNN</h2> <pre> from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n_neighbors=5, metric='minkowski', p=2) knn.fit(X_train, y_train) KNeighborsClassifier() y_pred = knn.predict(X_test) + Code + Markdown #Calculate accuracy of the model from sklearn.metrics import accuracy_score accuracy_KNN = accuracy_score(y_test, y_pred) print("Accuracy_KNN: {accuracy_KNN*100}") Accuracy_KNN: 96.17486338797814 </pre>

Final Model Selection Justification (2 Marks):

Final Model	Reasoning
Logistic Regression	C : Controls the regularization strength. Smaller values specify stronger regularization. Solver : Algorithm to use for optimization.
SVM	C : Controls the trade-off between achieving a low training error and a low testing error. Kernel : Defines the type of kernel function. Gamma : Determines the influence of a single training example.
Decision Tree	Max Depth : Limits the depth of the tree to prevent overfitting. Min Samples Split : Ensures that nodes are split only if a minimum number of samples is met.
Random Forest	n_estimators : The number of trees in the forest. Max Features : The number of features to consider when looking for the best split
KNN	n_neighbors : The number of neighbors to use for classification. Metric : The distance metric used for finding neighbors.