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Loading and Preprocessing

```
import pandas as pd
from sklearn.datasets import load_breast_cancer
from sklearn.preprocessing import StandardScaler

#loading the dataset
data = load_breast_cancer()
df = pd.DataFrame(data.data, columns=data.feature_names)
df['target'] = data.target

df.head()
```

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	mean fractal dimension	...	worst texture	worst perimeter	worst area	worst smoothness	worst compactness	worst concavity	worst concave points	worst symmetry	worst fractal dimension	target
0	17.99	10.38	122.80	1001.0	0.11840	0.27760	0.3001	0.14710	0.2419	0.07871	...	17.33	184.60	2019.0	0.1622	0.6656	0.7119	0.2654	0.4601	0.11890	0
1	20.57	17.77	132.90	1326.0	0.08474	0.07864	0.0869	0.07017	0.1812	0.05667	...	23.41	158.80	1956.0	0.1238	0.1866	0.2416	0.1860	0.2750	0.08902	0
2	19.69	21.25	130.00	1203.0	0.10960	0.15990	0.1974	0.12790	0.2069	0.05999	...	25.53	152.50	1709.0	0.1444	0.4245	0.4504	0.2430	0.3613	0.08758	0
3	11.42	20.38	77.58	386.1	0.14250	0.26390	0.2414	0.10520	0.2597	0.09744	...	26.50	98.87	567.7	0.2098	0.8663	0.6869	0.2575	0.6638	0.17300	0
4	20.29	14.34	135.10	1297.0	0.10030	0.13280	0.1980	0.10430	0.1809	0.05883	...	16.67	152.20	1575.0	0.1374	0.2050	0.4000	0.1625	0.2364	0.07678	0

5 rows x 31 columns

```
[2]: #checking missing values
df.isnull().sum()
```

	0
mean radius	0
mean texture	0
mean perimeter	0
mean area	0

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✓ 0s #checking missing values
df.isnull().sum()

	0
mean radius	0
mean texture	0
mean perimeter	0
mean area	0
mean smoothness	0
mean compactness	0
mean concavity	0
mean concave points	0
mean symmetry	0
mean fractal dimension	0
radius error	0
texture error	0
perimeter error	0
area error	0
smoothness error	0
compactness error	0
concavity error	0
concave points error	0
symmetry error	0
fractal dimension error	0
worst radius	0

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[2]

worst radius0

worst texture0

worst perimeter0

worst area0

worst smoothness0

worst compactness0

worst concavity0

worst concave points0

worst symmetry0

worst fractal dimension0

target0

1s

[3]

#feature scaling

scaler = StandardScaler()

scaled_features = scaler.fit_transform(df.drop('target', axis=1))

#DataFrame with scaled features

scaled_df = pd.DataFrame(scaled_features, columns=data.feature_names)

scaled_df['target'] = data.target

scaled_df.head()

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	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	mean fractal dimension	...	worst texture	worst perimeter	worst area	worst smoothness	worst compactness	worst concavity	worst concave points	worst symmetry	worst fractal dimension	target
0	1.097064	-2.073335	1.269934	0.984375	1.568466	3.283515	2.652874	2.532475	2.217515	2.255747	...	-1.359293	2.303601	2.001237	1.307686	2.616665	2.109526	2.296076	2.750622	1.937015	0
1	1.829821	-0.353632	1.685955	1.908708	-0.826962	-0.487072	-0.023846	0.548144	0.001392	-0.868652	...	-0.369203	1.535126	1.890489	-0.375612	-0.430444	-0.146749	1.087084	-0.243890	0.281190	0
2	1.579888	0.456187	1.566503	1.558884	0.942210	1.052926	1.363478	2.037231	0.939685	-0.398008	...	-0.023974	1.347475	1.456285	0.527407	1.082932	0.854974	1.955000	1.152255	0.201391	0
3	-0.768909	0.253732	-0.592687	-0.764464	3.283553	3.402909	1.915897	1.451707	2.867383	4.910919	...	0.133984	-0.249939	-0.550021	3.394275	3.893397	1.989588	2.175786	6.046041	4.935010	0
4	1.750297	-1.151816	1.776573	1.826229	0.280372	0.539340	1.371011	1.428493	-0.009560	-0.562450	...	-1.466770	1.338539	1.220724	0.220556	-0.313395	0.613179	0.729259	-0.868353	-0.397100	0

```
[2] worst symmetry 0
      worst fractal dimension 0
      target 0

dtype: int64
```

```
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5 rows x 31 columns

- Explanation:
- Dataset doesn't have missing values, missing value handling avoids potential errors during model training.
 - Used StandardScaler to scale the features to work better when the features are scaled to a similar range.

Classification Algorithm Implementation

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[3] #DataFrame with scaled features
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Explanation:

- Dataset doesn't have missing values, missing value handling avoids potential errors during model training.
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Classification Algorithm Implementation

[8] from sklearn.linear_model import LogisticRegression
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier
from sklearn.svm import SVC
from sklearn.neighbors import KNeighborsClassifier
from sklearn.model_selection import train_test_split
from sklearn.metrics import accuracy_score

#split the data
X = scaled_df.drop('target', axis=1)
y = scaled_df['target']

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classification_problem.ipynb ☆

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Classification Algorithm Implementation

```
[8] from sklearn.linear_model import LogisticRegression
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from sklearn.svm import SVC
from sklearn.neighbors import KNeighborsClassifier
from sklearn.model_selection import train_test_split
from sklearn.metrics import accuracy_score

#split the data
X = scaled_df.drop('target', axis=1)
y = scaled_df['target']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

#initialize models
models = {
    "logistic Regression" : LogisticRegression(),
    "Decision Tree" : DecisionTreeClassifier(),
    "Random Forest" : RandomForestClassifier(),
    "SVM" : SVC(),
    "k-NN" : KNeighborsClassifier()
}

#train and evaluate models
results = {}
for name, model in models.items():
    model.fit(X_train, y_train)
    y_pred = model.predict(X_test)
    acc = accuracy_score(y_test, y_pred)
    results[name] = acc
    print(f"{name} : Accuracy = {acc:.2F}")
```

Logistic Regression : Accuracy = 0.97
Decision Tree : Accuracy = 0.94
Random Forest : Accuracy = 0.96
SVM : Accuracy = 0.97
k-NN : Accuracy = 0.95

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[8] #train and evaluate models
results = {}
for name, model in models.items():
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 acc = accuracy_score(y_test, y_pred)
 results[name] = acc
 print(f"{name} : Accuracy = {acc:.2f}")

Logistic Regression : Accuracy = 0.97
Decision Tree : Accuracy = 0.94
Random Forest : Accuracy = 0.96
SVM : Accuracy = 0.97
k-NN : Accuracy = 0.95

Descriptions:

1. Logistic Regression : A linear model for binary classification that predict probabilities.
2. Decision Tree : A tree-based model that splits data based on feature values. It handles non-linear relationships effectively.
3. Random Forest : An ensemble of decision trees that reduces overfitting and improves generalization.
4. SVM : A model that finds a hyperplane to separate classes with maximum margin. Effective in high-dimensional spaces.
5. k-NN : A distance-based model that assigns a class based on the majority class among the nearest neighbors.

Model Comparison

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best_model = max(results, key=results.get)
worst_model = min(results, key=results.get)

print(f"The best performing model is {best_model} with an accuracy of {results[best_model]:.2f}")
print(f"The worst performing model is {worst_model} with an accuracy of {results[worst_model]:.2f}")

The best performing model is logistic Regression with an accuracy of 0.97
The worst performing model is Decision Tree with an accuracy of 0.94

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