Classification of 5G base stations

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# 1. Introduction

This project addresses the classification of such base stations using machine learning techniques applied to channel frequency response (CFR) data.

The objective is to build a classifier capable of identifying three classes:

* Class 0: Legitimate eNodeB (5G base station)
* Class 1: Attacker at location 1
* Class 2: Attacker at location 2

Using CFR samples and supervised learning, we aim to distinguish between these three classes with high accuracy (above 99%).

# 2. Problem Description

Each CFR sample is a 2D matrix with dimensions 72 (subcarriers) by 48 (repetitions). The training dataset contains labeled examples, while the test set is used for final evaluation through Kaggle. The core challenge is correctly classifying new, unseen data into one of the three predefined classes.

We have also noticed that the dataset is imbalanced, which means that class 0 is the dominant (around 80-90%) part of the dataset. Therefore, we decided to include a data augmentation block, particularly focusing on restoring the correct label distribution of the input data.

# 3. Data preprocessing

The whole architecture of our code is presented in the block diagram in **Fig.1**.

We started by loading all training ".npy" files based on the filenames listed in "label\_train.csv". Then we normalized the data using Z-score normalization - this helps stabilize training by ensuring the input has zero mean and unit variance.

Next, we split the dataset into training and validation sets using the "train\_test\_split". Here, we used a classical 80 to 20 split, which was actively used during the PC labs this semester.

A diagram of a process

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Fig. 1 Complete Block diagram of the presented solution

At this point, we understood that the dataset was imbalanced - class 0 (legit base stations) was much more frequent than classes 1 and 2 (attackers). To fix this, we implemented a custom data augmentation function. It adds Gaussian noise to underrepresented classes, especially classes 1 and 2, so all classes end up with similar counts. This was done before splitting the data to ensure the train and validation data have the same distribution of classes. The final distribution after augmentation and data splitting is presented in **Fig.2**.

A comparison of a bar graph

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Fig. 2 Class distribution in the training and validation sets after augmentation

# 4. Model architecture

To solve the task, we used a classical 2D Convolutional Neural Network (CNN). The model contains:

* One convolutional layer (Conv2D) with tunable filters and kernel size
* An optinal 2nd and 3rd convolutional layer (controlled via boolean variables)
* MaxPooling2D after each convolutional layer
* Flatten, 2 Dense layers (with tunable size)
* Dropout and final Sense layer with softmax activation

We also added a custom callback "EarlyStoppingWithThreshold", which works as a classical "EarlyStopping", but starts to monitor the required metrics only after it reaches a pre-defined threshold. The motivation behind it is that in some cases (mostly for complex models), it takes up to 10 epochs before the model starts to train (so that the "val\_accuracy" increases and "val\_loss" decreases). The custom callback was developed to wait over this period and prevent the training from stopping at this stage. One more callback we used is "ReduceLROnPlateau", which decreases the learning rate dynamically, if needed, and can help to achieve a high accuracy faster (in less epochs).

Therefore, the model is very flexible in terms of options, and by using a hyperparameter tuner, it is possible to find the simplest model out of the best-performing models.

To get the best configuration, we used Keras Tuner with "RandomSearch" option. Here, the tuner searches through a pre-defined number of random combinations (at the end, we set it to 10 combinations as a compromise between execution time and dispersion of models). The parameters we tuned include:

* Number of filters: 64, 128, 256, 512
* Kernel size: 3 or 5
* Size of the dense layer: from 64 to 512
* Dropout rate: from 0.2 to 0.5
* Learning rate: between 0.0001 and 0.01
* Boolean variables to include 2nd and 3rd convolutional layers

From the top 5 models with the highest validation accuracy, we picked the one with the smallest number of trainable parameters, as seen in **Fig.3**. The further training metrics for the best model are presented in **Fig.4**.

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Fig. 3 The best model picked out of five top-performing

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Fig. 4 Accuracy and loss graphs for the selected model

# 5. Results evaluation

After training and tuning, our model achieved 99.8% accuracy. The detailed results with misclassified detection are presented in the Confusion matrix in **Fig.5**. After predicting labels for test data and submitting it to Kaggle, the actual score is 0.99166, which is the top score at the moment and shows that the designed model is capable of predicting labels correctly in almost all cases.

A blue squares with numbers and labels

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Fig. 5 Confusion matrix of the best model

# 6. Conclusion

# 7. References

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