

Classifying EEG Spectrograms by Phalangeal Articulations utilizing Long-term Recurrent Convolutional Neural Networks

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1. Introduction

This design document contains information on the operation, architecture, API, and performance of the ESPA system, developed for the research on classifying EEG spectrograms by phalangeal articulations utilizing long-term recurrent convolutional (LRC) neural networks.

2. Theory of Operation

2.1. Data

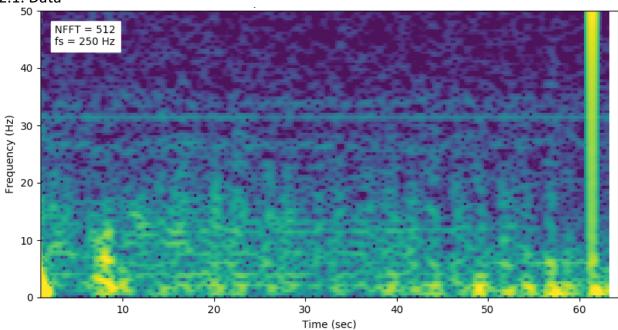


Figure 1 Sample spectrogram from 1 channel for sustained left middle finger flexion

For the inputs, the raw data consists of EEG time signals from 8 channels. The data from each channel is cleaned by performing the following steps:

- 1. Trim setup and teardown data
- 2. Remove DC offset
- 3. Notch mains interference
- 4. Bandpass filter frequencies from 1 to 50 Hz

After cleaning the data, spectrograms are computed, then partitioned into multiple samples with a dimensionality of 250 frequency points by 50 time points. Also, depending on the training run configuration, the samples are either replicated or augmented to increase the sample size. Finally, each sample is labelled with a one-hot encoded representation of its class. For example, [1., 0., 0.] would indicate class 1 in a 3-class model.

For the outputs, arrays containing predicted probabilities for each class are utilized, where the probabilities are represented in fractional form. For example, [0.75, 0.15, 0.10] would indicate 75% probability for class 1, 15% probability for class 2, and 10% probability for class 3.

2.2. Model

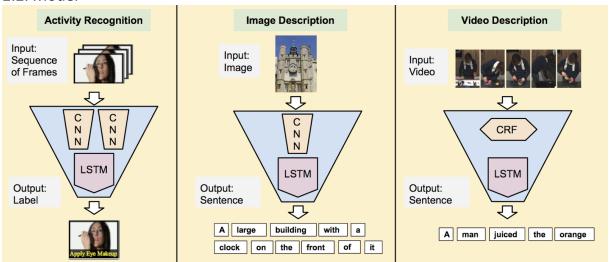


Figure 2 Applications of LRC neural networks (Source: http://jeffdonahue.com/lrcn/)

The model is based on long-term recurrent convolutional (LRC) neural networks, a class of neural networks used for visual and sequence learning ^[1]. It consists of a hybrid architecture of convolutional neural networks, recurrent neural networks, and multilayer perceptrons.

2.2.1. Convolutional Neural Networks (CNNs)

CNNs are biologically-inspired artificial neural networks that mimic the visual cortex. In a visual cortex, there are complex arrangements of cells that are sensitive to stimuli within a restricted region known as a receptive field. This region is tiled across an entire visual field, where the cells act as localized filters for detecting spatial patterns, the response to which can be approximated by a convolution operation [2][3]:

$$(f * g)[n] \stackrel{\text{def}}{=} \sum_{m=-\infty}^{\infty} f[m]g[n-m]$$
$$= \sum_{m=-\infty}^{\infty} f[n-m]g[m]$$

To illustrate how CNNs work, a sample 5x5 image, its grayscale conversion, and its simplified digital representation, where 1 is the maximum value instead of 255, is shown below:

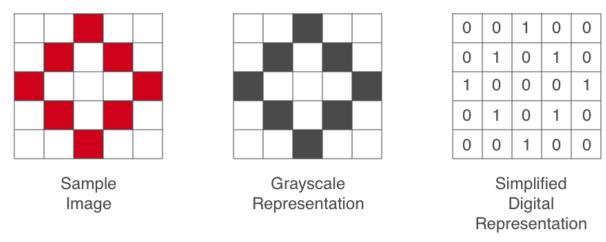


Figure 3 A sample image, its grayscale representation, and its simplified digital representation

A sample 3x3 filter and its simplified digital representation is also shown below:

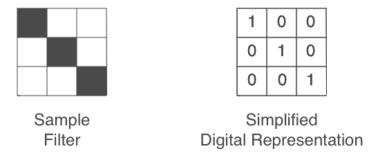


Figure 4 A sample filter and its simplified digital representation

In CNNs, convolutional filters are tiled across an image. As the filters tile across an image, receptive fields are convolved with their corresponding visual field regions, generating an activation map. In these activation maps, regions with with a high correlation with the filter pattern have high activation values, and vice versa. In this example, the filter has a stride of 1 pixel, generating a 3x3 activation map:

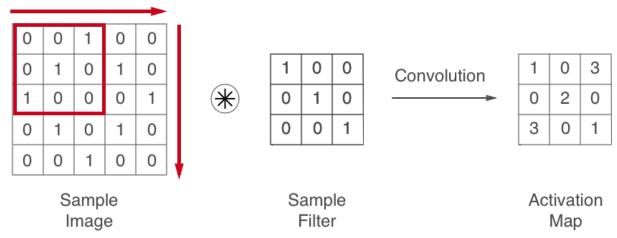


Figure 5 Generation of an activation map by tiling a filter across an image and performing successive convolutions

These activation maps are then passed through a layer of rectified linear units (ReLUs), an activation function used to improve the network's nonlinearity:

$$f(x) = \max(0, x)$$

Finally, a pooling layer downsamples the activation maps, reducing the number of parameters. In this example, a type of pooling layer called MaxPool is used, which replaces a pool of values with its maximum values, with a pool size of 2x2 and a stride of 1 pixel:

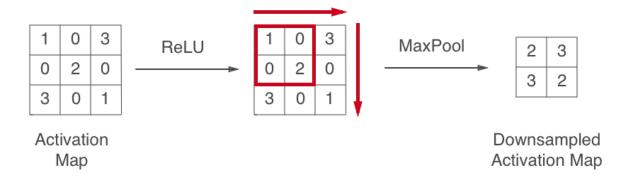


Figure 6 Application of a ReLU activation layer and a MaxPool pooling layer to an activation map

In practical applications, multiple alternating layers of convolution, ReLU activation, and MaxPool pooling are utilized. For this model, an architecture called VGG-16 [4] is used:

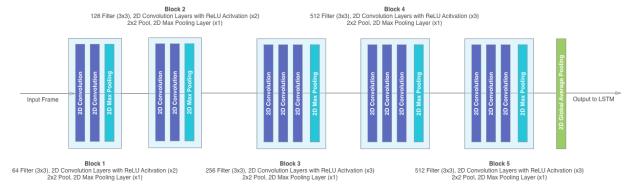


Figure 7 General VGG-16 architecture, adapted for this model

2.2.2. Recurrent Neural Networks (RNNs)

RNNs are artificial neural networks that are used for detecting sequential patterns. They consist of stateful memory units that are cyclically connected. One specific type of RNN, called long short-term memory (LSTM) ^[5] is used in the model. LSTMs, similar to regular RNNs, consist of chains of repeated LSTM units. However, unlike regular RNNs, LSTMs are well suited for data with variable gaps between events, such as variations observed in speech due to demographic and biological variability. To demonstrate how LSTMs work, diagrams and descriptions adapted from Colah's blog ^[6] and DeepLearning tutorials ^[7] are shown below:

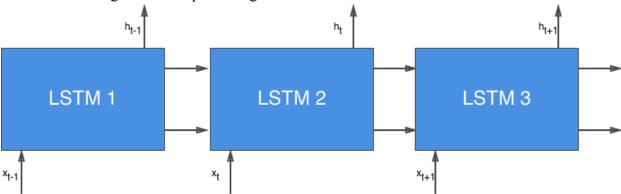


Figure 8 Chain of repeating LSTM units

Within each LSTM unit, several operations occur, which are represented by yellow circles on the diagram below:

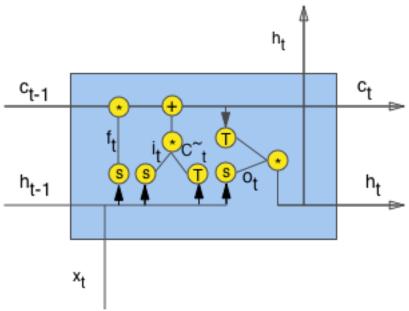


Figure 9 LSTM operations

In the diagram above, S represents a logistic sigmoid operation:

$$S(t) = \frac{1}{1 + e^{-t}}$$

T represents a hyperbolic tangent operation:

$$T(t) = \frac{e^{2t} - 1}{e^{2t} + 1}$$

+ represents element-wise addition, and * represents element-wise multiplication. LSTM operations also utilize weight matrices W, U, and V, and bias vector \mathbf{b} . First, the LSTM unit selects new data to store, which involves a logistic sigmoid layer (input gate) that selects which values to update:

$$i_t = S(W_i x_t + U_i h_{t-1} + b_i)$$

and a hyperbolic tangent layer that generates new candidate values:

$$C_t^{\sim} = T(W_c x_t + U_c h_{t-1} + b_c)$$

Next, the LSTM unit selects data to forget, which involves another logistic sigmoid layer:

$$f_t = S(W_f x_t + U_f h_{t-1} + b_f)$$

The layer takes in the input, x_t , and the previous output, h_{t-1} , then returns either 0 or 1 for each value in the cell state C_{t-1} , where 0 represents "forget" and 1 represents "remember". Then, the LSTM unit updates the cell state from the old state C_{t-1} to the new state C_t :

$$C_t = i_t C_t^{\sim} + f_t C_{t-1}$$

This operation forgets what has to be forgotten by multiplying the old cell state C_{t-1} with the output of the forget gate f_t , and adds new candidate values scaled by update weights by multiplying the new cell state C_t with the output of the input gate i_t . Finally, the LSTM unit generates the output. First, a logistic sigmoid layer selects which values of the cell state to output:

$$o_t = S(W_o x_t + U_o h_{t-1} + V_o C_t + b_o)$$

Next, the cell state values pass through a hyperbolic tangent layer, scaling the values between -1 and 1. Finally, the outputs are multiplied, resulting in a filtered cell state:

$$h_t = o_t T(C_t)$$

2.2.3. Multilayer Perceptrons (MLPs)

MLPs are artificial neural networks that consist of fully-connected layers of nodes. They map input data into outputs via a learned nonlinear transformation, which projects input data into a space where they become linearly separable, enabling classification:

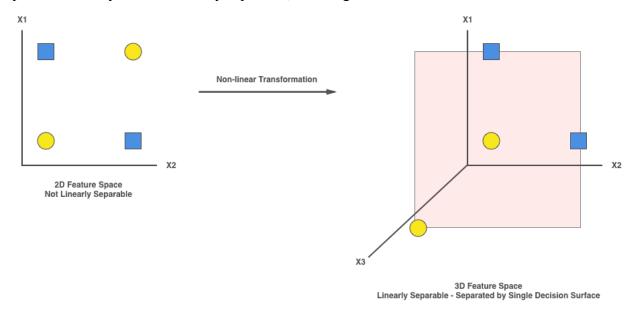


Figure 10 Simple illustration of how projecting input data into a feature space enables classification. Two classes are represented by a blue square and a yellow circle. A decision surface is represented by a red square.

MLPs consist of 3 primary stages: an input layer, hidden layers, and an output layer. With at least 1 hidden layer, an MLP becomes a universal approximator ^[8]. However, in practical deep learning applications, multiple hidden layers are utilized to generate more features. In the example below, an MLP with a 2-node input layer, 3-node hidden layer, and 2-node output layer is shown:

Hidden Layer

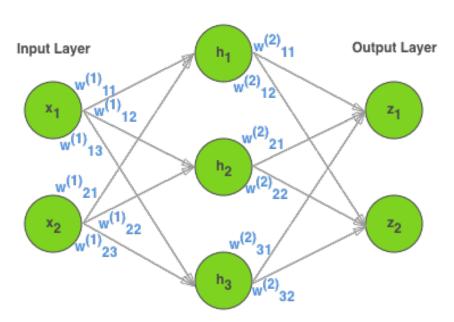


Figure 11 Sample MLP with a 2-node input layer, 3-node hidden layer, and 2-node output layer.

In MLPs, input nodes represent input features, hidden nodes represent generated features, and output nodes represent predicted class probabilities. To make predictions, an algorithm called forward propagation is used [8]:

$$h(x) = s(b^{(1)} + w^{(1)}x)$$

$$z(h(x)) = G(b^{(2)} + w^{(2)}h(x))$$

Where **x** is the input layer vector, **h** is the hidden layer vector, **z** is the output layer vector, **b** are bias vectors, **w** are weight matrices, **s** is the hidden layer activation function, which is set to ReLU for this model, and **G** is the output layer activation function, which is set to the softmax function for multi-class classification:

$$\sigma(z)_j = \frac{e^{z_j}}{\sum_{k=1}^K e^{z_k}}, j = 1, ..., K$$

Initially, the learned parameters from \mathbf{w} are randomized, resulting high error and low accuracy values. To improve accuracies, parameters are learned via the backpropagation algorithm [9][10], which trains the model on labelled data and updates parameters until a cost function is minimized.

3. System Architecture

3.1. Overview

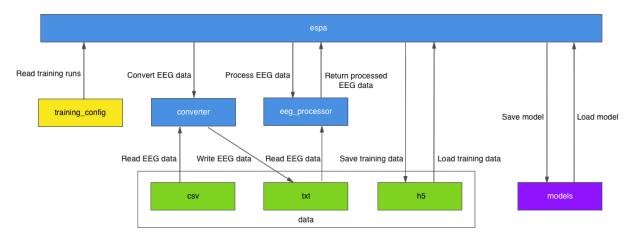


Figure 12 General system architecture

The diagram above shows the general system architecture of the ESPA system. Blue components represent Python modules, yellow components represent configuration files, green components represent data files, and purple components represent model files.

The espa module acts as the core module with which the user interfaces for executing training workflows. The converter module performs data type conversion operations for the raw EEG data. The eeg_processor module performs preprocessing operations on the raw EEG data. The training_config file enables the user to specify configurations for multiple training runs. The csv and txt data files store raw EEG data, while h5 data files store preprocessed training data. Finally, models store multiple model data, which come in pairs of model and weights data, stored in JSON and HDF5 files, respectively.

Multilayer Perceptron (1024 Hidden Units, 3 Output Units)

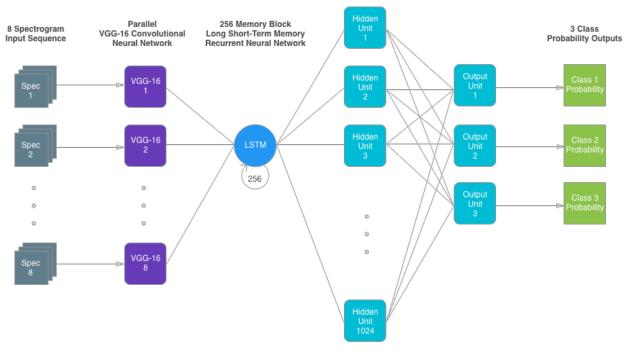


Figure 13 ESPA model architecture

The diagram above shows the ESPA model architecture. The model starts of with the input, which consists of a sequence of 8 spectrograms, each of which are replicated into 3 channels to match the expected input dimensions of the VGG-16 CNNs. Next, 8 VGG-16 CNNs processes the inputs in parallel. Then, a 256-memory block LSTM RNN processes the output of the previous stage. Finally, an MLP with a 1024-hidden unit hidden layer and 3-unit output layer processes the output of the previous layer and yields 3 class probability outputs.

4. Requirements

The ESPA system depends on the following:

- 1. Python, for general purpose programming
- 2. OpenCV, for computer vision
- 3. Keras, for implementing neural networks
- 4. Theano, for scientific and mathematical computations
- 5. SciPy, for scientific and mathematical computations
- 6. (Recommended) Nvidia GPU, for running computations in parallel

5. Application Programming Interface

5.1. Training Configuration File

The training configuration file, training_config.json, contains specifications for training runs. The configuration file consists of a list of training run specification, each of which have the following format:

```
"run_name": string name of training run,
    "trials": integer number of trials,
    "data_save_fn": string sample data save file path,
    "validation_ratio": float fraction of data allocated for validation,
    "testing_ratio": float fraction of data allocated for testing,
    "samples_generated_per_sample": integer number of samples generated/sample,
    "augmentation": boolean augmentation flag,
    "augmentation_magnitude": float augmentation magnitude,
    "freq_points": integer number of discrete spectrogram frequency points,
    "time_points": integer number of discrete spectrogram time points,
    "espa_save_fn": string ESPA model save file path,
    "espa_weights_save_fn": string ESPA model weights save file path
}
```

5.2. ESPA Module

The ESPA module, espa.py, provides the core user interface for executing training workflows.

5.2.1. class ESPAModel

Core class for instantiating and interfacing with ESPA models

5.2.1.1. method constructor

Constructor for the ESPAModel class.

Arguments:

- 1. data_save_fn: string sample data save file path
- 2. validation ratio: *float* fraction of data allocated for validation
- 3. testing_ratio: *float* fraction of data allocated for testing
- 4. samples generated per sample: integer number of samples generated/sample
- 5. augmentation: boolean augmentation flag
- 6. augmentation magnitude: float augmentation magnitude
- 7. freq_points: integer number of discrete spectrogram frequency points
- 8. time_points: integer number of discrete spectrogram time points
- 9. espa_save_fn: string ESPA model save file path
- 10. espa_weights_save_fn: string ESPA model weights save file path

Returns:

None

5.2.1.2. *method* train_espa_model

Trains the ESPA model

Arguments:

None

Returns:

1. metrics_history: Keras History object history of training and validation metric values

5.2.1.3. *method* test_espa_model

Tests the ESPA model

Arguments:

None

Returns:

1. metrics: dictionary testing categorical accuracy and loss

5.2.1.4. generator method generate_data

Generates data from HDF5 sample data file on demand

Arguments:

- 1. data_save_file: HDF5 file object sample data file
- 2. data_set: string data set to read
- 3. sample idxs: Numpy array array of sample indices
- 4. batch_size: integer number of samples for each batch

Yields:

1. Tuple of Numpy arrays containing a batch of samples and their labels

5.2.1.5. *method* print espa summary

Prints a summary representation of the ESPA model

Arguments:

None

Returns:

None

5.2.1.6. *method* generate_espa_model

Compiles the ESPA model

Arguments:

None

Returns:

None

5.2.1.7. *method* save_espa_model Saves the ESPA model and weights

Arguments:

None

Returns:

None

5.2.1.8. *method* load_espa_model Loads the ESPA model and weights

Arguments:

None

Returns:

None

5.2.1.9. *method* process_data Preprocesses sample data

Arguments:

None

Returns:

None

5.2.1.10. *method* replicate_augment_data Replicates or augments sample data

Arguments:

- 1. X_h: Numpy array pre replication or augmentation sample data
- 2. Y_h: *Numpy array* pre replication or augmentation sample data labels Returns:
 - 1. X: Numpy array post replication or augmentation sample data
 - 2. Y: Numpy array post replication or augmentation sample data labels
- 5.2.2. *class* ProgressDisplay

Callback class for displaying progress updates

5.2.2.1. *method* on batch end

Displays metric values at the end of each batch

Arguments:

- 1. epoch: *integer* epoch number
- 2. logs: dictionary log of metrics and their values

Returns:

None

5.2.3. Auxiliary Functions

5.2.3.1. function get training configuration

Acquires training configuration from a training configuration file Arguments:

- 1. training_config_fn: *string* training configuration file path Returns:
 - 1. training_config: list training configuration data

5.2.3.2. *function* execute_training_runs

Executes training runs from a specified training configuration

Arguments:

1. training config: list training configuration data

Returns:

1. results: dictionary compiled results data from training runs

5.2.3.3. function save results

Saves compiled results data to a JSON file

Arguments:

1. results_save_fn: string compiled results data save file path

Returns:

None

5.3. EEG Processor Module

The EEG processor module, eeg_processor.py, performs preprocessing operations on raw EEG data. This module was adapted from EEGrunt.py of the EEGrunt package, developed by Curiositry. For more information, please visit the EEGrunt repository: https://github.com/curiositry/EEGrunt.

5.4. Converter Module

The converter module, converter.py, performs data type conversion operations for raw EEG data. This module was adapted from convert_txt_to_csv.py of the EEGrunt package, developed by Curiositry. For more information, please visit the EEGrunt repository: https://github.com/curiositry/EEGrunt.

6. Training

6.1. Setup

The raw data comprises of 9, 8-channel EEG data, saved in text format, 3 for each of the following classes:

- 1. left index finger flexion
- 2. left middle finger flexion
- 3. left ring finger flexion

The data is then converted from text files into CSV files. Then, the data is filtered and trimmed to remove DC offset, mains interference, and setup/teardown artifacts. Next, spectrograms are calculated for each channel, each of which split into samples with 250 discrete frequency points and 50 discrete time points, and replicated 3 times to match the CNN's input dimensions, which expects 3-colour channel RGB inputs, generating a 30 x 8 x 3 x250 x 50 training dataset. Finally, the data is saved into an HDF5 file, which is vital for on-demand loading of data as a workaround for memory resource limitations.

The following training setups were implemented, with 3 trials per setup, and 10 epochs per trial:

- 1. Replication, 10x sample count
- 2. 1% augmentation, 10x sample count
- 3. 5% augmentation, 10x sample count

For each setup, the data was split into the following components:

- 1. 60% training data
- 2. 20% validation data
- 3. 20% testing data

6.2. Results and Discussion

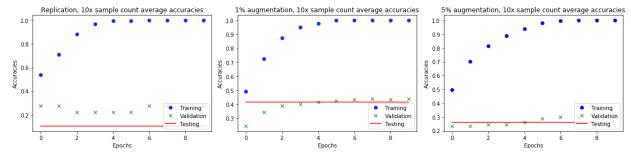


Figure 14 Accuracies (average training, validation, and testing) across all three setups

Across all three setups, it can be observed that the training accuracies ascend to high values and plateau within several epochs, while the validation accuracies lag behind the training accuracies, which could indicate overfitting. It can also be observed that both the validation and testing accuracies are significantly higher with 1% augmentation, which could indicate that 1% augmentation provides a good balance between generating independent samples and introducing excessive noise.

7. Recommendations

The primary issue that has to be addressed is overfitting, which prohibits the model from generalizing to new data. Some potential solutions include:

- 1. Increasing the raw sample size instead of depending entirely on data augmentation
- 2. Implementing regularization: L1, L2, and max norm
- 3. Implementing dropout

Other alternative changes that could potentially improve the model's performance include the choice of CNN architecture (e.g. ResNet, Inception), RNN architecture (e.g. Gated Recurrent Unit [GRU]), MLP architecture (deeper [more hidden layers] and wider [more neurons]). With an improved model, classification could be expanded to classify finer phalangeal articulations, given sufficient training data.

8. References

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- [9] C. V. D. Malsburg, "Frank Rosenblatt: Principles of Neurodynamics: Perceptrons and the Theory of Brain Mechanisms," Brain Theory, pp. 245–248, 1986.
- [10] D. Rumelhart, G. Hinton, and R. Williams, "Learning Internal Representations by Error Propagation," Readings in Cognitive Science, pp. 399–421, 1988.

9. Appendix

All source files can be found at the ESPA repository (https://github.com/valencra/eeg-spectrogram-phalangeal-articulation)

9.1. training_config.json

```
[ {
      "run name": "Replication, 10x sample count",
      "trials": 3,
      "data save fn": "data/h5/rep 10x data.h5",
      "validation ratio": 0.2,
      "testing ratio": 0.2,
      "samples generated per sample": 10,
      "augmentation": false,
      "augmentation magnitude": 0.00,
      "freq points": 250,
      "time points": 50,
      "espa save fn": "models/rep 10x espa model.json",
    "espa weights save fn": "models/rep 10x espa weights.h5"
}, {
      "run name": "1% augmentation, 10x sample count",
      "trials": 3,
      "data save fn": "data/h5/1pcaug 10x data.h5",
      "validation ratio": 0.2,
      "testing ratio": 0.2,
      "samples generated per sample": 10,
      "augmentation": true,
      "augmentation magnitude": 0.01,
      "freq points": 250,
      "time points": 50,
    "espa save fn": "models/1pcaug 10x espa model.json",
    "espa weights save fn": "models/1pcaug 10x espa weights.h5"
}, {
      "run name": "5% augmentation, 10x sample count",
      "trials": 3,
      "data save fn": "data/h5/5pcaug 10x data.h5",
      "validation ratio": 0.2,
      "testing ratio": 0.2,
      "samples generated per sample": 10,
```

```
"augmentation": true,
    "augmentation_magnitude": 0.05,
    "freq_points": 250,
    "time_points": 50,

"espa_save_fn": "models/5pcaug_10x_espa_model.json",
    "espa_weights_save_fn": "models/5pcaug_10x_espa_weights.h5"
}]
```

9.2. espa.py

```
from converter import convert txt to csv
from copy import deepcopy
from eeg processor import EEGProcessor
from keras import backend as K
from keras.applications.vgg16 import VGG16
from keras.callbacks import Callback
from keras.models import Model, model from json
from keras.layers import Dense, Dropout, Flatten, Input
from keras.layers.pooling import GlobalAveragePooling2D
from keras.layers.recurrent import LSTM
from keras.layers.wrappers import TimeDistributed
from keras.optimizers import Nadam
from keras.preprocessing.image import random rotation,
random shift, random shear, random zoom
from keras.utils.io utils import HDF5Matrix
from pprint import pprint
from math import ceil
from numpy import log10
from os import listdir
from os.path import join, isfile
from json import dump, load
import h5py
import numpy as np
import os
import sys
class ESPAModel(object):
      def init (self,
```

```
data save fn, validation ratio,
testing ratio, samples generated per sample,
                      augmentation, augmentation magnitude,
freq points, time points,
                      espa save fn, espa weights save fn):
             K.set image dim ordering("th")
             self.data save fn = data save fn
             self.validation ratio = validation ratio
             self.testing ratio = testing ratio
             self.samples generated per sample =
samples generated per sample
             self.augmentation = augmentation
             self.augmentation magnitude = augmentation magnitude
             self.freq points = freq points
             self.time points = time points
             self.espa save fn = espa save fn
             self.espa weights save fn = espa weights save fn
             self.espa = None
      def train espa model(self):
             """ Train the ESPA model
            print "\nTraining ESPA Model"
            batch size = 32
             with h5py.File(self.data save fn, "r") as
data save file:
                   # indices
                   training sample idxs =
np.random.permutation(range(int(data save file.attrs["training s
ample count"])))
                   validation sample idxs =
np.random.permutation(range(int(data save file.attrs["validation
sample count"])))
                   # generators
                   training sequence generator =
self.generate data(data save file=data save file,
```

```
data set="train",
sample_idxs=training_sample_idxs,
                                 batch size=batch size)
                   validation sequence generator =
self.generate data(data save file=data save file,
data set="val",
sample idxs=validation sample idxs,
                                   batch size=batch size)
                   # fit model
                   progress display = ProgressDisplay()
                   metrics history =
self.espa.fit generator(generator=training sequence generator,
validation data=validation sequence generator,
samples per epoch=len(training sample idxs),
nb val samples=len(validation sample idxs),
nb epoch=10,
verbose=2,
callbacks=[progress display],
class weight=None,
nb worker=1)
             return metrics history
```

```
def test espa model(self):
             """ Test the ESPA model
             print "\nTesting ESPA Model"
             batch size = 32
             with h5py.File(self.data save fn, "r") as
data save file:
                   # indices
                   testing sample idxs =
np.random.permutation(range(int(data save file.attrs["testing sa
mple count"])))
                   # generators
                   testing sequence generator =
self.generate data(data save file=data save file,
data set="test",
                                sample idxs=testing sample idxs,
                                batch size=batch size)
                   # calculate steps
                   sample count = len(testing sample idxs)
                   batches = int(sample count/batch size)
                   remainder samples = sample count%batch size
                   if remainder samples:
                          batches = batches + 1
                   # test model
                   metrics =
self.espa.evaluate generator(testing sequence generator,
batches)
                   # map metric names to metric values
```

```
metrics = {metric name: metric value for
metric name, metric value in zip(self.espa.metrics names,
metrics) }
                   print "Accuracy: {0:>8.4f} | Loss:
{1:>8.4f}".format(float(metrics["categorical accuracy"]),
                                        float (metrics["loss"]))
             return metrics
      def generate data(self, data save file, data set,
sample idxs, batch size):
             """ Generates data from HDF5 file on demand
             while True:
                   # determine batches
                   sample count = len(sample idxs)
                   batches = int(sample count/batch size)
                   remainder samples = sample count%batch size
                   if remainder samples:
                          batches = batches + 1
                   # generate batches
                   for idx in xrange(batches):
                          # incomplete batches
                          if idx == batches - 1:
                                batch idxs =
sample idxs[idx*batch size:]
                          # complete batches
                          else:
                                batch idxs =
sample_idxs[idx*batch size:idx*batch size+batch size]
                          batch idxs = sorted(batch idxs)
```

```
X = data save file[" ".join(["x",
data set])][batch idxs]
                          Y = data save file[" ".join(["y",
data set])][batch idxs]
                          yield (np.array(X), np.array(Y))
      def print espa summary(self):
             """ Prints a summary representation of the OSR model
             print "\n*** MODEL SUMMARY ***"
             self.espa.summary()
      def generate espa model(self):
             """ Builds the ESPA model
             print "\nGenerating ESPA model..."
             with h5py.File(self.data save fn, "r") as
data save file:
                   class count =
len(data save file.attrs["classes"].split(","))
             # input layer
             spectrograms = Input(shape=(8,
                                                        3,
self.freq points,
self.time points))
             # CNN layers
             cnn base = VGG16(input shape=(3,
self.freq points,
self.time points),
                                        weights="imagenet",
                                        include top=False)
```

```
cnn out = GlobalAveragePooling2D()(cnn base.output)
             cnn = Model(input=cnn base.input, output=cnn out)
             cnn.trainable = False
             encoded spectrograms =
TimeDistributed(cnn) (spectrograms)
             # RNN layers
             encoded spectrograms =
LSTM(256) (encoded spectrograms)
             # MLP layers
             hidden layer = Dense(output dim=1024,
activation="relu") (encoded spectrograms)
             outputs = Dense(output dim=class count,
activation="softmax") (hidden layer)
             # compile model
             espa = Model([spectrograms], outputs)
             optimizer = Nadam(lr=0.0002,
                                         beta 1=0.9,
                                         beta 2=0.999,
                                         epsilon=1e-08,
                                         schedule decay=0.004)
             espa.compile(loss="categorical crossentropy",
                                 optimizer=optimizer,
metrics=["categorical accuracy"])
             self.espa = espa
      def save espa model(self):
             """ Save the ESPA model to an HDF5 file
             # delete save files, if they already exist
             try:
                   print "\nESPA save file \"{0}\" already
exists! Overwriting previous saved
file.".format(self.espa save fn)
                   os.remove(self.espa save fn)
```

```
except OSError:
                   pass
             try:
                   print "ESPA weights save file \"{0}\" already
exists! Overwriting previous saved
file.\n".format(self.espa weights save fn)
                   os.remove(self.espa weights save fn)
             except OSError:
                   pass
             # save ESPA model
            print "\nSaving ESPA model to
\"{0}\"".format(self.espa save fn)
            with open(self.espa save fn, "w") as espa save file:
                   espa model json = self.espa.to json()
                   espa save file.write(espa model json)
             # save ESPA model weights
            print "Saving ESPA model weights to
\"{0}\"".format(self.espa weights save fn)
             self.espa.save weights (self.espa weights save fn)
            print "Saved ESPA model and weights to disk\n"
      def load espa model(self):
             """ Load the ESPA model from an HDF5 file
            print "\nLoading ESPA model from
\"{0}\"".format(self.espa save fn)
            with open(self.espa save fn, "r") as espa save file:
                   espa model json = espa save file.read()
                   self.espa = model from json(espa model json)
            print "Loading ESPA model weights from
\"{0}\"".format(self.espa weights save fn)
            with open(self.espa weights save fn, "r") as
espa weights save file:
```

```
self.espa.load weights (self.espa weights save fn)
            print "Loaded ESPA model and weights from disk\n"
      def process data(self):
             """ Preprocesses data
            print "\nProcessing data..."
             data dirs = sorted(["left-index-flexion",
                                             "left-middle-
flexion",
                                             "left-ring-
flexion"])
            txt data dir = "data/txt"
             csv data dir = "data/csv"
             # convert text data into CSV data
             for data dir in data dirs:
                   convert txt to csv(join(txt data dir,
data dir),
join(csv data dir, data dir))
            X = {channel idx:[] for channel idx in range(8)}
             Y = []
            data sample count = 0
             # iterate through all class directories
             for class idx, data dir in enumerate(data dirs):
                   class dir = join(csv data dir, data dir)
                   class files = [class file
                                   for class file in
listdir(class dir)
                                   if (isfile(join(class dir,
class_file))) and (".csv" in class_file)]
```

```
sys.stdout = open(os.devnull, "w") # silence
EEG data processing standard outputs
                   # iterate through all class files
                   for class file in class files:
                          session title = "
".join(class file.split("-"))
                          eeg processor =
EEGProcessor(class dir, class file, "openbci", session title)
                          eeg processor.plot = 'show'
                          eeg processor.load data()
                          # iterate through all channels
                          for channel idx, channel in
enumerate(eeg processor.channels):
                                print " ".join(["Processing
channel ",
      str(channel idx + 1))
                                # load and clean channel data
      eeg processor.load channel(channel)
                                eeg processor.remove dc offset()
      eeg processor.notch mains interference()
                                eeg processor.trim data(10, 10)
                                # calculate spectrogram
                                eeg processor.get spectrum data()
                                eeg processor.data =
eeg processor.bandpass(1, 50)
                                spec =
10*log10(eeg processor.spec PSDperBin)
                                # accumulate sample data
```

```
sample count = spec.shape[1] /
self.time points
                                for sample idx in
xrange(sample count):
                                       sample =
spec[0:self.freq points,
sample idx*self.time points:
sample idx*self.time points+self.time points]
                                       format spec = lambda
spectrogram: np.array([spectrogram] *3)
      X[channel idx].append(format spec(sample))
                          # accumulate label data
                          data sample count += sample count
                          label = [0] * len (data dirs)
                          label[class idx] = 1
                          label = np.array(label)
                          Y.extend([label] * (sample count))
                   sys.stdout = sys. stdout # stop silencing
standard outputs
             # format sample and label data
             X = [np.array([cha 1, cha 2, cha 3, cha 4, cha 5,
cha 6, cha 7, cha 8])
                    for cha 1, cha 2, cha 3, cha 4, cha 5,
cha 6, cha 7, cha 8
                        in zip(X[0], X[1], X[2], X[3],
                                X[4], X[5], X[6], X[7])
             X = np.array(X)
             Y = np.array(Y)
             # save sample and label data into HDF5 file
```

```
with h5py.File(self.data save fn, "w") as
data save file:
                    # partition data into training, validation,
and testing sets
                   sample idxs =
np.random.permutation(range(data sample count))
                   training sample idxs =
sample idxs[0:int((1.0-self.validation ratio-
self.testing ratio) *data sample count)]
                   validation sample idxs =
sample idxs[int((1.0-self.validation ratio-
self.testing ratio) *data sample count):int((1.0-
self.testing ratio) *data sample count)]
                    testing sample idxs = sample idxs[int((1.0-
self.testing ratio) *data sample count):]
                   x train, y train =
self.replicate augment data(X[training sample idxs],
Y[training sample idxs])
                   x val, y val =
self.replicate augment data(X[validation sample idxs],
Y[validation sample idxs])
                   x \text{ test}, y \text{ test} =
self.replicate augment data(X[testing sample idxs],
Y[testing sample idxs])
                   data save file.attrs["classes"] =
np.string (",".join(data dirs))
                    data save file.attrs["training sample count"]
= len(x train)
      data save file.attrs["validation sample count"] =
len(x val)
                   data save file.attrs["testing sample count"]
= len(x test)
                    # training set
```

```
x train ds =
data save file.create dataset("x train",
shape=(len(x train), 8, 3, self.freq points, self.time points),
dtype="f")
                   y train ds =
data save file.create dataset("y train",
shape=(len(y train), len(data dirs)),
dtype="i")
                   x train ds[:] = x train
                    y train ds[:] = y train
                    # validation set
                    x val ds =
data save file.create dataset("x val",
shape=(len(x val), 8, 3, self.freq points, self.time points),
dtype="f")
                    y val ds =
data save file.create dataset("y val",
shape=(len(y val), len(data_dirs)),
dtype="i")
                    x \text{ val } ds[:] = x \text{ val}
                    y_val_ds[:] = y_val
                    # testing set
                    x test ds =
data save file.create dataset("x test",
shape=(len(x test), 8, 3, self.freq points, self.time points),
dtype="f")
```

```
y test ds =
data save file.create dataset("y_test",
shape=(len(y test), len(data dirs)),
dtype="i")
                    x \text{ test } ds[:] = x \text{ test}
                    y \text{ test } ds[:] = y \text{ test}
       def replicate augment data(self, X h, Y h):
              """ Replicates/augments sample data
              11 11 11
             # samples after replication/augmentation
             X = []
             Y = []
             # spectrogram formatting function
             format spec = lambda spectrogram:
np.array([spectrogram] *3)
             # replicate/augment samples
             for x h, y h in zip(X h, Y h):
                    # increase sample data via image augmentation
                    if self.augmentation:
                           for in
xrange(self.samples generated per sample):
                                  x = [] # new sample
                                  for spectrogram in x h:
                                         spectrogram =
spectrogram[0] # first colour channel only
                                         shifted sample =
random shift(np.array([spectrogram]),
wrg = self.augmentation magnitude,
hrg = self.augmentation magnitude)
```

```
x.append(format spec(shifted_sample[0]))
                                X.append(np.array(x))
                                Y.append(y h)
                   # increase sample data via image replication
                   else:
                          for in
xrange(self.samples generated per sample):
                                X.append(x h)
                                Y.append(y h)
             # replicated/augmented samples
            return np.array(X), np.array(Y)
class ProgressDisplay(Callback):
      """ Progress display callback
      def on batch end(self, epoch, logs={}):
            print " Batch {0:<4d} => Accuracy: {1:>8.4f} |
Loss: {2:>8.4f} | Size: {3:>4d}".format(int(logs["batch"])+1,
                    float(logs["categorical accuracy"]),
                    float(logs["loss"]),
                    int(logs["size"]))
# auxilliary functions
def get training configuration(training config fn):
      """ Acquires training configuration from a file
      with open(training config fn, "r") as
training config file:
```

```
training config = load(training config file)
      return training config
def execute training runs (training config):
      """ Executes training runs from specified training
configuration
      11 11 11
      results = {}
      # iterate through training runs
      for training run in training config:
             run name = training run["run name"]
             trials = training run["trials"]
             data save fn = training run["data save fn"]
             validation ratio = training run["validation ratio"]
             testing ratio = training run["testing ratio"]
             samples generated per sample =
training run["samples generated per sample"]
             augmentation = training run["augmentation"]
             augmentation magnitude =
training run["augmentation magnitude"]
             freq points = training run["freq points"]
             time points = training run["time points"]
             espa save fn = training run["espa save fn"]
             espa weights save fn =
training run["espa weights save fn"]
             results[run name] = {}
             print "\n".join(["="*80,
                                 "EXECUTING TRAINING RUN:
{0}".format(run name),
                                 "="*801)
             # iterate through trials
             for trial in xrange(trials):
                   print "\n".join(["-"*80,
                                       "TRIAL:
{0}".format(trial),
                                     "-"*80])
```

```
espa = ESPAModel(data save fn = data save fn,
                                              validation ratio =
validation ratio,
                                              testing ratio =
testing ratio,
samples generated per sample = samples generated per sample,
                                              augmentation =
augmentation,
augmentation magnitude = augmentation magnitude,
                                               freq points =
freq points,
                                              time points =
time points,
                                              espa save fn =
espa save fn,
                                            espa weights save fn
= espa weights save fn)
                   espa.process data()
                   espa.generate espa model()
                   espa.print espa summary()
                   training metrics history =
espa.train espa model()
                   testing metrics = espa.test_espa_model()
                   espa.save espa model()
                   # organize results
                   results[run name][trial] = {}
                   results[run name][trial]["train"] =
training metrics history.history
                   results[run name][trial]["test"] =
testing metrics
      return results
def save results(results, results save fn):
```