Abstract—A major issue for bringing brain-computer interface (BCI) based on electroencephalogram (EEG) recordings outside of laboratories is the non-stationarities of EEG signals. Varying statistical properties of the signals during inter- or intra-session transfers can lead to deteriorated BCI performances over time. These variations may cause the input data distribution to shift when transitioning from the training phase (calibration session) to the testing/operating phase resulting in a covariate shift. We propose to handle this issue using a novel hybrid learning method based on two classifiers, wherein the first classifier allows including new information in the training dataset, and the second classifier performs an overall classification. The proposed method is motivated by the smoothness assumption, i.e., the points that are closest to each other are more likely to share the same label, and may be added online to enrich the training dataset. The method is evaluated on two real-world datasets corresponding to motor imagery detection (BCI competition 2008 dataset 2A and 2B). The results support the conclusion that an improvement in the classification accuracy over traditional inductive learning and semi-supervised learning methods can be obtained.

I. INTRODUCTION

A Brain-Computer Interface (BCI) is an alternative communication’s means, which allows a user to express his will without muscle exertion, provided that the brain signals are properly translated into computer commands [1], [2], [3]. With an electroencephalography (EEG) based BCI that operates online in real-time non-stationary/changing environments, it is required to consider input features that are invariant to shifts of the data, or learning approaches that can be able to track the changes that can repeat overtime, to update the classifier in a timely fashion. In fact, it may be difficult to reliably classify the EEG patterns in BCI during long sessions using traditional inductive classification algorithms due to the non-stationarity characteristics of the EEG signal. The non-stationarities in the EEG may be caused by various reasons such as changing user attention level, electrode placement, and user fatigue [4], [5], [6]. There are therefore notable variations or shifts in the EEG signals during trial-to-trial, and session-to-session transfers [7]. These variations often appear as covariate shifts, wherein the input data distributions differ significantly between training/calibration and testing/operating phases, while the conditional distribution remains the same [8], [9].

The low classification accuracy has been one of the main concerns of the developed BCI systems based on motor imagery (MI) detection, which directly affects the reliability of the BCI decision making process and the information transfer rate. The traditional classification algorithms are mainly inductive. They deal with the development of a function that approximates output values using input data from the whole searching space (i.e. induction), and then uses this function to predict the output values for all the new input vectors (i.e. deduction). To enhance BCI performance, several feature extraction, feature selection, and feature classification techniques have been proposed in the literature [10], [11], [12], [13]. A large variety of features have been used in BCI based on motor imagery detection such as band powers, power spectral density, time frequency features, and common spatial patterns (CSP) based features [14]. However, the EEG non-stationary characteristics result in shifts in feature distributions for all types of features in varying degrees. An example showing the change of distribution between the training and test datasets is depicted in Figure 1.

To solve this issue, adaptive learning algorithms have been proposed for devising adaptive BCI systems with positive results [4], [15], [16]. Most of which have made efforts to reduce the non-stationarity aspect in the extracted features. In an adaptive learning technique, a priori information is required about the changes in the EEG signals. Additionally, the adaptive techniques are mostly based on supervised learning techniques, which need labeled data (i.e. a training dataset that can be obtained through a calibration session [7], [17]), which may not be available or difficult to acquire during the operating phase. Different strategies are possible, first the same classifier can be used over time, and the classifier is updated only when a significant shift is detected [18], [19]. Second, it can be assumed that the signal will always change over time in a continuous way, and not severe shift can be detected. In such a case, the classifier will be updated with all the new trials that can be correctly detected, and the updated training database will follow the data distribution. The later approach is chosen in the present study. It is also motivated by the time between two trials in motor imagery detection, which can be several seconds, and allows a classifier to be fully retrained thanks to the addition of new knowledge.
To overcome the issues discussed above, a transductive-inductive learning approach based on two classifiers is presented in this paper. The idea of the proposed learning algorithm is to handle the covariate shift by initiating the adaptation based on both the existing (training dataset) and new knowledge obtained through the testing phase. The transductive classifier is only used for adding new information in the existing training dataset, and the inductive classifier is used for predicting the BCI outputs, after being retrained each time the training dataset is updated. It is thus a learning approach wherein transductive and inductive learning are combined to update the training dataset, and to track the evolution of the features over time. The first classifier is a probabilistic weighted K-nearest neighbor (PWKNN) based transductive classifier. The output of the first classifier is used to determine if a trial and its corresponding estimated label can be added to the training dataset, and the inductive learning model is updated. Transductive learning combines induction and deduction in a single step, and is related to the field of semi-supervised learning (SSL), which uses both labeled and unlabeled data during learning [20]. By eliminating the need to construct a global model, transductive learning offers prospect to achieve higher accuracy. In order to make use of unlabeled data, it is necessary to assume some structure to the underlying distribution of data. Additionally, it is essential that the SSL approach must satisfy at least one of the following assumptions such as smoothness, cluster, or manifold assumption [21], [22], [23]. In SSL, the widely used type of algorithm is graph based label propagation, which propagates labels from labeled examples to the whole unlabeled data. In the present paper, we will make use of the smoothness assumption (i.e., the points which are closest to each other are more likely to share the same label) to implement a transductive learning algorithm. The second classifier is inductive, a linear support vector machine (SVM) classifier, its outputs are used to determine the BCI outputs. In this paper, we demonstrate the effectiveness of the approach on a synthetic dataset and on two real-world datasets, and we show that the covariate shifts can be tracked and adapted using the proposed method. Specifically, using the data from the BCI competition-IV 2A and 2B, we demonstrate the effectiveness of the proposed approach over traditional, adaptive, and semi-supervised learning algorithms. The contributions of the paper can be summarized as follows: First, a combination of a transductive and inductive classifier are used to track, and address the effect of covariate shifts in the non-stationary EEG signals. Second, the proposed learning strategy updates automatically the classifier online without making any a priori assumption about the distribution for the upcoming test data. The remainder of the paper is organized as follows: section II presents the system overview of the transductive-inductive learning method, including the probabilistic K-nearest neighbor. The datasets and the signal processing steps are described in section IV. The results are then presented in section V and discussed in section VI.

II. System overview

A. Problem Statement

Let us consider a learning framework in which a training dataset is denoted by $X_{Tr} = \{(x_i, y_i)\}_{i=1}^N$, where $N$ is the total number of observations, and a label $y_i$ is associated with each input $x_i$. Depending upon the number of inputs and outputs, $x_i$ and $y_i$ may be a scalar or vector variables. Let us consider a two-class classification problem i.e., $y \in \omega_1, \omega_2$. The probability distribution of the inputs at time $i$ can thus be defined as, $P(x_i) = P(\omega_1)P(x_i|\omega_1) + P(\omega_2)P(x_i|\omega_2)$ where $P(\omega_1)$ and $P(\omega_2)$ are the prior probabilities of getting a sample of the classes $\omega_1$ and $\omega_2$, respectively, while $P(x_i|\omega_1)$ and $P(x_i|\omega_2)$ are the conditional probability distribution for the time period $i$. The goal is to predict the labels $\hat{y}_i$ of upcoming samples resulting into $X_{Tr} = \{\hat{y}_i|x_i\}_{i=1}^M$, where $M$ is the total number of observations in the testing phase.

B. Probabilistic K-Nearest Neighbor

Probability theory plays a vital role in the solution of pattern recognition problems, because most of the problems can be solved using a density estimation technique [24]. Its main task is to model a probability density function $P(x)$ of a random variable $X$, given $X_{Tr}$. There are two approaches for the density estimation, namely, parametric and non-parametric. One of the key limitations of the parametric approach is that it assumes a precise practical form for the distribution, which may lead to be incompatible for a specific application. An alternative approach is the non-parametric density estimation, which estimates density function without applying any assumption about underlying data distribution. Here, we consider a non-parametric method based on K-nearest-neighbors (KNN) because it is a transductive learning method that uses the test data point to determine a decision [25]. In a KNN method, we consider a small sphere centered at the point $x$ at which, we wish to estimate the density $P(x)$. We allow the radius of the sphere to grow until it contains $K$ data points, the estimate of the density is then given by:

$$P(x) = K/(N \cdot V)$$  \hspace{1cm} (1)$$

where, $V$ is set to the volume of the sphere, and $N$ is the total number of points. The parameter $K$ governs the degree of smoothing. The technique of KNN density estimation may be extended to the classification task in which the KNN density estimation is obtained for each class, and then the Bayes theorem is used to perform a classification task. Now, let’s suppose that we have a data set comprising $N_\omega$ points in the class $\omega_i$ within the set of classes $\omega$, where $i \in 1, 2$, so that $\sum_{\omega_i} N_\omega_i = N$. If we wish to classify a new point $x$, we draw a sphere centered on $x$ containing precisely $K$ points irrespective of their classes. Now suppose this sphere has the volume $V$, and contains $K(\omega_i)$ from class $\omega_i$. Then, an estimate of the density associated with each class or
Fig. 1. Covariate shift in the EEG dataset 2A subject A03, between training and testing input distribution for different frequency bands. (a) Mu band [8-12] Hz, and (b) Beta band [14-30] Hz. The red circle denote the features of the left hand motor imagery and blue crosses denote the features of the right hand motor imagery. The black and red lines represent the decision boundaries obtained by the training data and test data respectively.

The likelihood can be obtained by:

$$P(x|\omega_i) = \frac{K_{\omega_i}}{N_i \cdot V}$$

Similarly, the unconditional density is given by

$$P(x) = \frac{K}{N \cdot V}$$

Using the Bayes theorem, we can obtain the posterior probability of the class membership:

$$P(\omega_i|x) = \frac{P(x|\omega_i)P(\omega_i)}{P(x)} = \frac{K_{\omega_i}}{K}$$

If we wish to minimize the probability of misclassification, this is achieved by assigning the test point $x$ to the class $\omega_i$ having the largest posterior probability, i.e. corresponding to the largest value of $K_{\omega_i}/K$. Thus, to classify a new point, identify the K-nearest points from the training data set, and then assign the new point to the set having the largest number of representatives. This posterior probability is also known as the Bayesian belief or confidence ratio ($CR$). However, the overall estimate obtained by the KNN method may not be satisfactory, because the resulting density is not a true probability density since its integral over all the samples space diverges [26]. Another drawback is that it considers only the $K$ points to build the density, and each neighbor has an equal weight. An extension to the above KNN method is to assign the weight to each sample that depends on the distance to $x$ [27]. A radial basis function (RBF) kernel is used to obtain the weights. Using RBF Kernel, the nearest points have weights with higher value than furthest points. A probabilistic weighted K-nearest neighbors (PWKNN) approach based on an RBF kernel is thus proposed to devise the transductive classifier with $RBF_{(p,q)}$:

$$RBF_{(p,q)} = \exp(-\frac{d^2(p,q)}{2\sigma^2})$$

where $d(p,q)$ is the Euclidean distance from the unlabeled data point $x_p$ to the labeled data point $x_q$ is computed as given below:

$$d_{(p,q)} = \sqrt{\sum_{i=1}^{m} (x_p(i) - x_q(i))^2}$$

and $x(i)$ is the i-th feature of $x$ and $m$ is the number of features.

For binary detection, the confidence ratio of $CR_{\omega_1}$ of the class $\omega_1$, for a data point $x_p$, is defined by:

$$CR_{\omega_1} = \frac{\sum_{j=1}^{K} RBF_{(p,j)} \cdot (l_j = \omega_1)}{\sum_{j=1}^{K} RBF_{(p,j)}}$$

$$CR_{\omega_2} = 1 - CR_{\omega_1}$$

where $x_j$, $1 \leq j \leq K$, corresponds to the j-th nearest neighbor of $x_p$. The outputs of PWKNN include the overall confidence of the decision, given by:

$$CR = \max(CR_{\omega_1}, CR_{\omega_2})$$

and the output class $\tilde{y}$ (1 if $x_p$ is assigned to $\omega_1$, 0 otherwise).

III. CLASSIFIER COMBINATION

In the proposed system, a transductive classifier is used to determine if an example should be used or not to enrich a training dataset, which is used by an inductive classifier to determine the final output. Transductive learning was proposed and briefly studied more than 40 years ago by Vapnik and Chervonenk [25]. Later, it has been empirically acknowledged that transduction can often serve as a more efficient, or accurate learning method than traditional inductive learning approaches [22]. This appreciation is the rationale behind the development of the transductive learning model...
for accounting non-stationarities in BCIs. Moreover, transductive learning methods estimate the value of a potential model (classification function) at a new data point using specific training cases related to that test point [25]. Using an SSL approach, every unlabeled point gives information about \( P(x) \). The usefulness of an example depends on the condition whether the density of the input distribution has changed significantly, or the point is located very close to the classification boundary. For example, in SSL, the smoothness assumption states that the points closest to each other are more likely to share the same label. Hence, if the point satisfies the smoothness assumption, then a single observation provides some useful information to update the training data-set. Then, using the updated training data-set, an improved global model can be obtained to classify the new unlabeled data points.

In this paper, we propose a novel transductive-inductive learning model. The proposed approach consists of two steps: induction and transduction. Initially, an inductive classifier, denoted by \( \mathcal{F} \), is trained on the features obtained from the calibration/training data \( X_{Tr} \). Once the classifier \( \mathcal{F} \) is trained, and a classification decision boundary is obtained, then the evaluation phase starts. In this phase, we set the parameters \( CR_{\alpha}, K, \Delta m \), where \( CR_{\alpha} \) is a confidence ratio threshold, \( K \) is the number of neighbors for preparing a transductive classifier (denoted by \( \mathcal{T} \)) through PWKNN algorithm, and \( \Delta m \) is the number of data points after which the adaptation starts. In addition, the classifier \( \mathcal{F} \) will classify the input features obtained from the testing data \( X_{Ts} \) in the evaluation phase. The classifier \( \mathcal{F} \) will initiate adaptation through transductive learning after every \( \Delta m \) consecutive trials (in this study, we have fixed \( \Delta m = 5 \)). Considering a small number of trials \( \Delta m \) in each epoch, results in focusing on a trial-by-trial shift, which may be caused by various reasons such as muscular artifacts. Each time the classifier \( \mathcal{F} \) initiates adaptation, it is considered as one epoch, and it assigns \( \Delta m \) data points from \( X_{Ts} \) to a variable \( \Delta X_{Ts} \), and predicts the labels through the transductive function \( \mathcal{T} \). In this step, each observation is taken from \( \Delta X_{Ts} \), and the transductive function \( \mathcal{T} \) is applied. Once all the \( \Delta X_{Ts} \) points are processed through \( \mathcal{T} \), the \( CR \) value is used for each trial to decide if the trial’s features and the corresponding estimated output should be added to \( X_{Tr} \), i.e. if \( CR > CR_{\alpha} \). Thus, the labels \( \mathcal{T}(\Delta X_{Ts}) \) obtained through \( \mathcal{T} \), which are above \( CR_{\alpha} \) are inserted into the training dataset. Based on the updated training dataset, the inductive function \( \mathcal{F} \) is updated (i.e. the classifier is retrained), and a new classifier is obtained. This process is repeated until all the \( M \) points in the testing phase are classified. The pseudo-code of the algorithm is given in the Algorithm 1.

**Algorithm 1 Online training and evaluation.**

1. Train \( \mathcal{F} \) with \( X_{Tr} \)
2. Set the parameters \( CR_{\alpha}, K, \Delta m \)
3. for each trial \( x_i \) from \( X_{Ts} \) do
4. \( \hat{y}_i = \mathcal{F}(x_i) \)
5. if \( i \mod \Delta m \equiv 0 \) then
6. \( \Delta X_{Ts} = X_{Ts}(i - m + 1 : i) \)
7. for each trial \( x_j \) from \( \Delta X_{Ts} \) do
8. \( \{ \mathcal{T} = PWKNN \} \)
9. \( (CR, \hat{y}) = PWKNN(X_{Tr}, x_j, CR_{\alpha}, K) \)
10. if \( CR > CR_{\alpha} \) then
11. Updated \( X_{Tr} = X_{Tr} + (x_j, \hat{y}) \)
12. Train \( \mathcal{F} \) with \( X_{Tr} \)

**B. Data Processing and Feature Extraction**

The first stage of signal processing employs a filter bank that decomposes the EEG signals into multiple frequency bands [29]. A total of 10 band-pass filters are used, namely [8-12], [10-14], [12-16], [14-18], [16-20], [18-22], [20-24], [22-26], [24-28], [26-30] Hz [30]. This set of frequency bands is used because it covers the expected frequency range of motor imagery response. In the next sections, we consider a time segment of 3 s after the cue onset for both data sets (BCI Competition IV dataset 2A and dataset 2B).

The second stage employs common spatial filters (CSP), a set of spatial filters that maximize the variance of spatially filtered signals under one condition, while minimizing it for the other condition. Raw EEG scalp potentials are known to have poor spatial resolution due to volume conduction. If the signal of interest is weak while other sources produce strong...
signals in the same frequency range, then it is difficult to classify two classes of EEG measurements [14].

C. Evaluation

In order to evaluate the performance of the system, we have considered the classification accuracy as the measure of performance index. A Cohens kappa coefficient (kappa value) is used to compare the performance with other competing methods. The experiments are performed using a linear SVM pattern classifier (F) and PWKNN classifier (T) for transductive learning. The accuracy is given in percentage (%). The hyper-parameters $K$ and $CR_{\alpha}$ needs to be carefully selected in order to limit the number of wrongly labeled elements that are added in the training database. Two variants of the proposed learning method, namely TI$_1$ and TI$_2$, are presented. In TI$_1$, the hyper-parameters are selected based on grid search to maximize the mean accuracy across subjects, with $K \in \{6, 12, 18\}$, and $CR_{\alpha} \in \{0.5, \ldots, 1\}$. In TI$_2$, the hyper-parameters are determined for each subject, they are selected to maximize the accuracy of each subject. In the dataset 2A, the session I is divided into two parts, 80% for training the classifier and 20% is used to determine the hyper-parameters. The evaluation is then performed on the data from the session II. In the dataset 2B, the sessions I and II are used for training $F$, session III is used to obtain the hyper-parameters, and sessions IV and V are used to evaluate the performance. For each dataset, the accuracy corresponding to a 10-fold-cross validation (10-CV) on the training dataset is presented. Additionally, two variants of the proposed method are presented and compared with the state-of-the-art methods. The baseline method uses a filter-bank with traditional inductive learning [29]. The baseline state-of-the-art methods. The baseline method uses a filter-training dataset is presented. Additionally, two variants of evaluating the performance. For each dataset, the accuracy of hyper-parameters, and sessions IV and V are used to classify two classes of EEG measurements [14].

The proposed learning method for EEG-based BCI is based on Vapnik’s principle of transduction i.e., given a test feature corresponding to a frequency band. The blue crosses and red circles denote the features of the left hand and right hand MI, respectively. The black line represents a possible linear separation between the features of the two classes obtained from each frequency band. The hyperplan is plotted for an illustration purpose only.

The performance results of the proposed learning algorithm are compared against the aforementioned methods, on the BCI Competition IV dataset 2A and dataset 2B. The accuracy for each subject, with the mean and standard deviation (std), is presented in Tables II and III. For the dataset 2A, with the baseline method, the accuracy is $77.78 \pm 14.63$ across subjects. The three variations of SSL provides an accuracy of $75.54 \pm 12.99$, $75.16 \pm 15.03$, and $75.95 \pm 14.73$ for SSL$_1$, SSL$_2$, and SSL$_3$. The proposed method provides an accuracy of $78.01 \pm 12.86$ for TI$_1$, and $79.17 \pm 13.3$ for TI$_2$. For the dataset 2B, the accuracy is $65.46 \pm 25.25$ with the baseline method, $65.50 \pm 13.63$, $65.59 \pm 13.63$, and $67.38 \pm 13.75$ for SSL$_1$, SSL$_2$, and SSL$_3$. Finally, TI$_1$ and TI$_2$ provide an accuracy of $69.56 \pm 26.4$ and $70.85 \pm 25.52$, respectively.

A two-sided Wilcoxon signed rank test is used to assess the statistical significance of the improvement at a confidence level of 0.05 in all the pairwise comparisons. The values of hyper-parameters $CR_{\alpha}$ and $K$ are presented in Fig. 3. For the dataset 2A, they are set as $CR_{\alpha} = 0.95$ and $K = 12$. For the dataset 2B, they are set as $CR_{\alpha} = 0.95$ and $K = 18$. For the dataset 2A, there is a margin of 0.23% in the average accuracy for the TI$_1$ method over the baseline method. Additionally, the proposed method has shown an improvement of 2.06% in average accuracy in comparison with SSL based label propagation methods. Moreover, the improvement over the baseline method is 4.6% for dataset 2B, but it is not statistically significant ($p=0.0938$). The values of $CR_{\alpha}$ and $K$ are given in Tables II and III for the dataset 2A and dataset 2B. The average accuracy for TI$_2$ is superior to the accuracy of the baseline method and the other reported methods for both the datasets 2A and 2B. However, for the dataset 2A, the improvement is only 1.39%, and it is not significant. However, there is statistically significant improvement of 5.39% ($p = 0.0391$) for the dataset 2B.

The effectiveness of proposed method for dataset 2A is presented in Table I, by comparing its classification results in terms of Kappa values. The TI$_1$ and TI$_2$ methods have better mean Kappa values over other competing methods. In addition, there is no statistically significant difference between TI$_1$ and TI$_2$ for both the databases 2A ($p = 0.25$) and 2B ($p = 0.21$). Finally, the average accuracy obtained with $UB$ is only 84.57% and 79.02% for the dataset 2A and 2B, respectively. It shows that the use of the test datasets with the expected ground truth does not allow to obtain an accuracy above 90% but the variability across sessions leads to a drop of 6.79% for 2A, and 13.56% for 2B.

VI. DISCUSSION

The proposed learning method for EEG-based BCI is based on Vapnik’s principle of transduction i.e., given a test
point, one should focus on the training points which are in a neighborhood of this test point. A PWKNN based on an RBF kernel is used to construct a local decision rule, and predict the label of the test point according to this transduction rule. The proposed approach has satisfied the smoothness assumption: If two instances are close to each other in a high-density region; they are likely to share the same label. The new information/knowledge obtained through transduction is used to update the training dataset of the inductive classifier. The main classification function is still inductive because the transductive knowledge is only used to add more information into training dataset.

A certain minimum value of confidence ratio ($CR_{\alpha}$) is used to decide whether the information is useful or not to enrich the classifier. If it is the case then it is added to the training dataset. The discarded information may come from artifacts or an abrupt change in the input distribution. The values of $CR_{\alpha}$ and $K$ needs to be carefully chosen in order to achieve better performance. With subject dependent hyperparameters values, the proposed TI$_2$ method shows a statistically significant improvement over the baseline method ($p < 0.05$).
TABLE II
CLASSIFICATION ACCURACY (% RESULTS FROM BCI COMPETITION IV DATASET 2A.

<table>
<thead>
<tr>
<th>Subject</th>
<th>10-CV Training</th>
<th>Baseline SSL</th>
<th>SSL$_1$ Eval</th>
<th>SSL$_2$ Eval</th>
<th>SSL$_3$ Eval</th>
<th>T$_1$ Eval</th>
<th>T$_2$ Eval</th>
<th>UB Eval</th>
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<td>A01</td>
<td>87.86</td>
<td>90.97</td>
<td>86.81</td>
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<td>A02</td>
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<td>61.11</td>
<td>56.94</td>
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TABLE III
CLASSIFICATION ACCURACY (% RESULTS FROM BCI COMPETITION IV DATASET 2B.

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<th>Subject</th>
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<th>SSL$_2$ Eval</th>
<th>SSL$_3$ Eval</th>
<th>T$_1$ Eval</th>
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<td>61.88</td>
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<td>76.56</td>
<td>71.88</td>
<td>71.88</td>
<td>75.94</td>
<td>89.69</td>
<td>18</td>
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</tr>
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<td>66.56</td>
<td>70.00</td>
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<td>64.06</td>
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<td>61.56</td>
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</tr>
<tr>
<td>B09</td>
<td>70.81</td>
<td>65.46</td>
<td>65.59</td>
<td>65.59</td>
<td>67.38</td>
<td>69.56</td>
<td>6</td>
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<tr>
<td>Mean</td>
<td>70.81</td>
<td>65.46</td>
<td>65.59</td>
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<td>67.38</td>
<td>69.56</td>
<td>70.85</td>
<td>79.02</td>
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<tr>
<td>Std</td>
<td>10.78</td>
<td>25.25</td>
<td>13.63</td>
<td>13.63</td>
<td>13.75</td>
<td>26.4</td>
<td>25.52</td>
<td>28.43</td>
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</tbody>
</table>

TABLE I
COMPARISON OVER TYPES OF EEG-BASED CLASSIFICATION METHODS ON KAPPA VALUES FROM BCI COMPETITION IV DATASET 2A.

<table>
<thead>
<tr>
<th>Subjects</th>
<th>CSP[32]</th>
<th>FBCSP (DC)[34]</th>
<th>RQNN[33]</th>
<th>T$_1$</th>
<th>T$_2$</th>
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<tbody>
<tr>
<td>A01</td>
<td>0.56</td>
<td>0.71</td>
<td>0.22</td>
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<td>0.88</td>
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<tr>
<td>A02</td>
<td>0.31</td>
<td>0.37</td>
<td>0.22</td>
<td>0.22</td>
<td>0.22</td>
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<tr>
<td>A03</td>
<td>0.70</td>
<td>0.66</td>
<td>0.58</td>
<td>0.76</td>
<td>0.88</td>
</tr>
<tr>
<td>A04</td>
<td>0.44</td>
<td>0.47</td>
<td>0.21</td>
<td>0.39</td>
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<tr>
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<td>0.22</td>
<td>0.41</td>
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<td>0.43</td>
<td>0.53</td>
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<td>A06</td>
<td>0.20</td>
<td>0.26</td>
<td>0.22</td>
<td>0.33</td>
<td>0.33</td>
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<tr>
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<td>0.61</td>
<td>0.73</td>
<td>0.17</td>
<td>0.38</td>
<td>0.38</td>
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<tr>
<td>A08</td>
<td>0.76</td>
<td>0.58</td>
<td>0.35</td>
<td>0.85</td>
<td>0.85</td>
</tr>
<tr>
<td>A09</td>
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<td>0.58</td>
<td>0.81</td>
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<tr>
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<td>0.33</td>
<td>0.56</td>
<td>0.58</td>
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<tr>
<td>Std</td>
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<td>0.16</td>
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VII. CONCLUSION

In this paper, we have proposed a hybrid classifier combination using transductive and inductive classifiers to address the effect of covariate shifts in non-stationary EEG signals associated with motor imagery detection in the EEG signal. The labels from the unlabeled data are estimated with a probabilistic weighted K-nearest neighbor approach. The proposed T$_2$ method that has subject dependent hyperparameters provides a statistically significant superior classification accuracy over a purely inductive baseline method on the dataset-2B. Particularly this learning approach provides a foundation for combining the transductive learning and inductive learning for EEG based BCI. Further work will include the detection of the covariate shift in the data, and only retrain the inductive classifier when a relevant shift is detected. This may reduce the processing time and complexity in re-training the classifier in the absence of any real change in the data.

VIII. ACKNOWLEDGMENTS

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Another important issue is the number of trials after which an adaptation is initiated. Considering a small number of trials in each epoch results in focusing on trial-by-trial shift, which can be due to muscular artifacts. However, the long term non-stationarities may be accounted for by considering a large number of trials in each epoch. We chose a small number of trials for updating the classifier, as our aim was to track the covariate shift trial-by-trial. The proposed transductive-inductive learning technique makes use of on-line data to extract features, and thus adapts to non-stationarities in the streaming EEG. The experimental results demonstrated the effectiveness of the proposed learning strategy, showing better performance than traditional learning methods.
REFERENCES


