
Online Multi-Task Learning based on K-SVD

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Abstract

This paper develops an efficient online algorithm based on K-SVD for learning multiple consecutive tasks. We first derive a batch multi-task learning method that builds upon the K-SVD algorithm, and then extend the batch algorithm to train models online in a lifelong learning setting. The resulting method has lower computational complexity than other current lifelong learning algorithms while maintaining nearly identical performance. Additionally, the proposed method offers an alternate formulation for lifelong learning that supports both task and feature similarity matrices.

1. Introduction

With the increasing interest in big data and deployed machine learning software, it is essential to develop systems that are capable of learning multiple, consecutive tasks over an extended time period. The idea of learning and sharing knowledge between multiple tasks to improve collective performance has been studied extensively (Kumar & Daumé III, 2012; Kang et al., 2011; Zhang et al., 2008; Caruana, 1997) from a batch learning perspective, in which all models are trained simultaneously. Recent work in online multi-task learning (MTL) (Ruvolo & Eaton, 2013; Saha et al., 2011) has shown that it is possible to learn tasks sequentially and achieve nearly identical accuracy to batch MTL methods while dramatically reducing the computational requirements. These capabilities are essential to the development of lifelong learning algorithms that continually accumulate and refine knowledge of multiple tasks over an unbounded stream of experience.

To facilitate knowledge transfer between task models, one common technique used by MTL algorithms

is to learn and maintain a shared repository of latent model components; each task model is then given as a weighted combination of these latent components. This technique is used successfully by several current MTL methods, including various batch algorithms (Kumar & Daumé III, 2012; Zhang et al., 2008) and the online Efficient Lifelong Learning Algorithm (ELLA) (Ruvolo & Eaton, 2013). In batch MTL algorithms, the latent model components are learned simultaneously with the task models in an expensive joint optimization. ELLA employs several simplifications to eliminate the expensive optimization in support of online learning while minimizing the adverse effect on the performance of the resulting task models. Most notably, ELLA requires that each model, once learned, be based on fixed weights over the latent components. This may adversely affect performance by not permitting task models to adjust the weighting of individual latent components as these components become more refined due to additional training.

In this paper, we investigate an alternate formulation of online multi-task learning based on the K-SVD algorithm (Aharon et al., 2006) that provides a reduction in computational complexity over ELLA, providing improved support for the rapid learning of consecutive tasks. This K-SVD formulation also eliminates one of ELLA's simplifications, enabling task models to flexibly adjust their weights over model components during learning. We compare this formulation to the original ELLA, a new version of ELLA that incorporates an iterative update step with a similar computational complexity to our K-SVD based algorithm, and a hybrid approach that blends these two approaches. We show that in some situations our K-SVD based algorithm exhibits similar learning performance to ELLA while allowing continual refinement of all models and latent components. We also show that the hybrid approach yields a robust algorithm that exhibits high performance across a range of learning domains.

We begin by providing an overview of the K-SVD algorithm, and then adapt K-SVD to learn task models in a batch MTL setting, yielding a new algorithm we

call MTL-SVD. We then modify the batch MTL-SVD algorithm to operate online, making it suitable for application to lifelong learning settings.

2. The K-SVD Algorithm

This section reviews the K-SVD algorithm of Aharon et al. (2006) for learning dictionaries for sparse coding, which forms the foundation of our approach. Suppose we are designing a dictionary consisting of k vectors to sparsely code a set of points $\{\mathbf{x}_1, \dots, \mathbf{x}_n\} \subset \mathbb{R}^d$. We would like to compute a dictionary $\mathbf{L} \in \mathbb{R}^{d \times k}$ such that each input point can be coded with a minimal number of dictionary elements. This objective can be realized by solving the following optimization problem:

$$\arg \min_{\mathbf{L}} \sum_{i=1}^n \min_{\mathbf{s}^{(i)}} \left\{ \|\mathbf{L}\mathbf{s}^{(i)} - \mathbf{x}_i\|_2^2 + \mu \|\mathbf{s}^{(i)}\|_0 \right\}, \quad (1)$$

where $\mathbf{s}^{(i)}$ is the vector of coefficients over the columns of \mathbf{L} to encode \mathbf{x}_i and μ is a positive constant that defines the tradeoff between accurate reconstruction of the input points and the sparsity of the coefficient vectors. This objective is computationally hard to optimize due to the cross terms between the dictionary \mathbf{L} and coefficients $\mathbf{S} = [\mathbf{s}^{(1)} \dots \mathbf{s}^{(n)}]$ as well as the presence of the L_0 norm $\|\cdot\|_0$, which both make the objective non-convex. Some approaches for solving Equation 1 alternately optimize \mathbf{L} and \mathbf{S} until a local minima is reached.¹

Like other approaches for dictionary learning, K-SVD alternates two optimization steps.

1. Optimize \mathbf{S} in Equation 1 given the current \mathbf{L} .
2. For a particular dictionary element (i.e., the j th column of \mathbf{L}), jointly optimize the element as well as its corresponding coefficient for each data point currently encoded by the dictionary element (i.e., the non-zero entries in the j th row of \mathbf{S}).

We next describe each of these steps of K-SVD; the complete K-SVD algorithm is given as Algorithm 1.

Step 1: Optimizing \mathbf{S}

Given a fixed value of \mathbf{L} , Equation 1 decomposes into n independent optimization problems of the form:

$$\mathbf{s}^{(i)} \leftarrow \arg \min_{\mathbf{s}} \left\{ \|\mathbf{L}\mathbf{s} - \mathbf{x}_i\|_2^2 + \mu \|\mathbf{s}\|_0 \right\}. \quad (2)$$

Equation 2 is known as the *sparse coding problem*, and can be solved (approximately) using numerous tech-

¹Optimizing \mathbf{L} given a fixed \mathbf{S} is a convex optimization problem, whereas optimizing the columns of \mathbf{S} with fixed \mathbf{L} , while not convex, can be relaxed into a convex optimization problem by replacing the L_0 norm with the L_1 norm.

Algorithm 1 K-SVD (Aharon et al., 2006)

input data points $\{\mathbf{x}_1, \dots, \mathbf{x}_n\}$, dictionary size k
 init \mathbf{L} using random column vectors of unit length
loop until convergence **do**
 for $i \in \{1, \dots, n\}$, perform update in Eqn. 2
 for $j \in \{1, \dots, k\}$, perform updates in Eqns. 4–6
end loop
return \mathbf{L}

niques (e.g., Matching Pursuit, Orthogonal Matching Pursuit, or the Lasso (Tibshirani, 1996)).

Step 2: Optimizing a Dictionary Element and its Corresponding Non-Zero Coefficients

This step updates a particular dictionary element as well as the corresponding coefficients for data points that are encoded using the element (i.e., have a non-zero coefficient value). Let \mathbf{l}_j indicate the particular column of \mathbf{L} to optimize. First, we form the matrix \mathbf{E} representing the residual for each data point given that \mathbf{l}_j is zeroed out. The i th column of \mathbf{E} is given by:

$$\mathbf{e}_i = \mathbf{x}_i - \sum_{r \neq j} s_r^{(i)} \mathbf{l}_r, \quad (3)$$

where $s_r^{(i)}$ is the r th entry of $\mathbf{s}^{(i)}$. Next, we perform a singular value decomposition (SVD) on \mathbf{E} . The first left singular vector provides the updated value for \mathbf{l}_j and the corresponding right singular vector scaled by the corresponding singular value yields the updated coefficients for each data point (i.e., the j th row of \mathbf{S}).

We would like both steps (1) and (2) to either maintain or improve the quality of our solution to Equation 1. Unfortunately, using the SVD of \mathbf{E} will cause some coefficients in \mathbf{S} that were previously zero to become non-zero, eliminating the guarantee that the quality of our solution cannot become worse. To eliminate this possibility, we take the SVD of the subset \mathcal{A} of the columns of \mathbf{E} such $m \in \mathcal{A} \Leftrightarrow \mathbf{s}_j^{(m)} \neq 0$:

$$(\mathbf{U}, \mathbf{\Sigma}, \mathbf{V}) = \text{svd}(\mathbf{E}_{\mathcal{A}}) \quad (4)$$

$$\mathbf{l}_j \leftarrow \mathbf{u}_1 \quad (5)$$

$$\mathbf{s}_j^{(\mathcal{A})} \leftarrow \sigma_{1,1} \mathbf{v}_1, \quad (6)$$

where $\mathbf{E}_{\mathcal{A}}$ denotes the matrix formed from the subset of columns in \mathcal{A} , the singular values are assumed to all be positive (this is possible for any real matrix) and sorted in descending order, and $\mathbf{s}_j^{(\mathcal{A})}$ denotes the vector formed from the columns in \mathcal{A} of the j th row of \mathbf{S} . It is well-known that this optimization procedure minimizes $\|\mathbf{E}_{\mathcal{A}} - \mathbf{B}\|_2^2$ for all rank-1 matrices $\mathbf{B} = \mathbf{l}_j \mathbf{s}_j^{(\mathcal{A})\top}$.

This implies that the resulting \mathbf{L}_j and $\mathbf{s}_j^{(A)}$ also minimize $\sum_{i=1}^n \|\mathbf{L}\mathbf{s}^{(i)} - \mathbf{x}_i\|_2^2$, which shows that quality of the solution to Equation 1 cannot have worsened.

3. Multi-Task Learning Using K-SVD

This section extends the K-SVD algorithm to batch multi-task learning. The key step of our approach is to adapt the K-SVD algorithm from the objective of learning a dictionary for sparse coding a set of input data points to learning a dictionary for sparse coding a set of parameter vectors for individual task models. We begin by describing the multi-task learning setting, and then describe how to adapt K-SVD to this setting. We call our resulting algorithm *MTL-SVD*.

3.1. Problem Setting

In the batch multi-task learning setting, the agent simultaneously learns models for a set of supervised learning tasks $\{\mathcal{Z}^{(1)}, \mathcal{Z}^{(2)}, \dots, \mathcal{Z}^{(T)}\}$. Each learning task $\mathcal{Z}^{(t)} = (\hat{f}^{(t)}, \mathbf{X}^{(t)}, \mathbf{y}^{(t)})$ is defined by a (hidden) mapping $\hat{f}^{(t)} : \mathcal{X}^{(t)} \mapsto \mathcal{Y}^{(t)}$ from an instance space $\mathcal{X}^{(t)} \subseteq \mathbb{R}^d$ to a set of labels $\mathcal{Y}^{(t)}$ (typically $\mathcal{Y}^{(t)} = \{-1, +1\}$ for classification tasks and $\mathcal{Y}^{(t)} = \mathbb{R}$ for regression tasks). Task t has n_t training instances $\mathbf{X}^{(t)} \in \mathbb{R}^{d \times n_t}$ with corresponding labels $\mathbf{y}^{(t)} \in \mathcal{Y}^{(t)n_t}$ given by $\hat{f}^{(t)}$. We assume that the learner is given all training data (instances and labels) for all learning tasks in a single batch. The agent's goal is to construct task models $f^{(1)}, \dots, f^{(T)}$ where each $f^{(t)} : \mathbb{R}^d \mapsto \mathcal{Y}^{(t)}$ such that each $f^{(t)}$ will approximate $\hat{f}^{(t)}$ to enable the accurate prediction of labels for new instances.

3.2. Model of Task Structure

Our model of latent task structure is based on the GO-MTL model proposed by Kumar & Daumé III (2012), which allows for learning of overlap and grouping between tasks. Kumar & Daumé III showed that their model performs well on a range of multi-task problems.

We assume a parametric framework in which each task model $f^{(t)}(\mathbf{x}) = f(\mathbf{x}, \boldsymbol{\theta}^{(t)})$ is specified by a task-specific parameter vector $\boldsymbol{\theta}^{(t)} \in \mathbb{R}^d$. To facilitate transfer between tasks, we maintain a library of k latent model components $\mathbf{L} \in \mathbb{R}^{d \times k}$ that are shared between task models. Each task parameter vector $\boldsymbol{\theta}^{(t)}$ can be represented as a linear combination of the columns of \mathbf{L} according to the coefficient vector $\mathbf{s}^{(t)} \in \mathbb{R}^k$ (i.e., $\boldsymbol{\theta}^{(t)} = \mathbf{L}\mathbf{s}^{(t)}$). We encourage the $\mathbf{s}^{(t)}$'s to be sparse (i.e., use few latent components) to ensure that each learned latent model component captures a maximal reusable chunk of knowledge.

Given the labeled training data for each task, we optimize the task models to minimize the predictive loss over all tasks while encouraging the models to share structure through \mathbf{L} . This optimization problem is realized by the objective function:

$$e_T(\mathbf{L}) = \sum_{t=1}^T \min_{\mathbf{s}^{(t)}} \left\{ \frac{1}{n_t} \sum_{i=1}^{n_t} \mathcal{L} \left(f \left(\mathbf{x}_i^{(t)}; \mathbf{L}\mathbf{s}^{(t)} \right), y_i^{(t)} \right) + \lambda \|\mathbf{L}\mathbf{s}^{(t)}\|_2^2 + \mu \|\mathbf{s}^{(t)}\|_0 \right\}, \quad (7)$$

where $(\mathbf{x}_i^{(t)}, y_i^{(t)})$ is the i th labeled training instance for task t , \mathcal{L} is a known loss function for fitting the task models, and λ and μ are non-negative constants that define the amount of ridge-regression on the models and the penalty for non-sparse solutions, respectively.

In earlier work (Ruvolo & Eaton, 2013), we proposed a method for optimizing Equation 7 efficiently. In part, we achieve this by recasting the objective function in Equation 7 as a problem of learning a dictionary for sparse-coding task models learned on the training data for each task in isolation. We accomplish this by taking the second-order Taylor expansion of Equation 7 about $\mathbf{L}\mathbf{s}^{(t)} = \boldsymbol{\theta}^{(t)}$, where $\boldsymbol{\theta}^{(t)}$ is an optimal predictor for task t learned on only that task's training data. We apply a similar technique to arrive at a simplified form of the original objective:

$$g_T(\mathbf{L}) = \sum_{t=1}^T \min_{\mathbf{s}^{(t)}} \left\{ \frac{1}{n_t} \|\boldsymbol{\theta}^{(t)} - \mathbf{L}\mathbf{s}^{(t)}\|_{\mathbf{D}^{(t)}}^2 + \mu \|\mathbf{s}^{(t)}\|_0 \right\} \quad (8)$$

where

$$\mathbf{D}^{(t)} = \lambda \mathbf{I} + \nabla_{\boldsymbol{\theta}, \boldsymbol{\theta}}^2 \frac{1}{2n_t} \sum_{i=1}^{n_t} \mathcal{L} \left(f \left(\mathbf{x}_i^{(t)}; \boldsymbol{\theta} \right), y_i^{(t)} \right) \Big|_{\boldsymbol{\theta}=\boldsymbol{\theta}^{(t)}}$$

$$\boldsymbol{\theta}^{(t)} = \arg \min_{\boldsymbol{\theta}} \frac{1}{n_t} \sum_{i=1}^{n_t} \mathcal{L} \left(f \left(\mathbf{x}_i^{(t)}; \boldsymbol{\theta} \right), y_i^{(t)} \right) + \lambda \|\boldsymbol{\theta}\|_2^2,$$

and $\|\mathbf{v}\|_{\mathbf{A}}^2 = \mathbf{v}^\top \mathbf{A} \mathbf{v}$. Note that Equation 8 is identical to Equation 1 except that we use the norm $\|\cdot\|_{\mathbf{D}^{(t)}}$ as opposed to $\|\cdot\|_2^2$. Therefore, we could ignore the difference in the objective functions and apply the unmodified K-SVD algorithm to the $\boldsymbol{\theta}^{(t)}$'s to arrive at a sensible multi-task learning algorithm. However, our goal will be to improve upon this naïve approach.

Our principal modification to the original K-SVD algorithm is to replace the SVD in Equation 4 with a generalized SVD:²

$$(\mathbf{U}, \boldsymbol{\Sigma}, \mathbf{V}) = \text{gsvd}(\mathbf{E}_A, \mathbf{M}, \mathbf{W}), \quad (9)$$

²Note that computing the GSVD is not more computationally expensive than computing the SVD, with the exception that we must first compute the matrix square roots of \mathbf{M} and \mathbf{W} .

where \mathbf{M} is a symmetric positive semidefinite (PSD) matrix in $\mathbb{R}^{d \times d}$ and \mathbf{W} is a symmetric PSD matrix in $\mathbb{R}^{T \times T}$. Later we will discuss how to choose \mathbf{M} and \mathbf{W} ; however, for now we assume they are given. In this case, the GSVD step defined in Equation 9 along with the updates in Equations 5 and 6 yield the values of \mathbf{l}_j and $\mathbf{s}_j^{(A)}$ that minimize:

$$\sum_{t_1=1}^{|\mathcal{A}|} \sum_{t_2=1}^{|\mathcal{A}|} w_{t_1, t_2} \left(\mathbf{e}_{\mathcal{A}_{t_1}} - \mathbf{l}_j \mathbf{s}_j^{(A_{t_1})} \right)^\top \mathbf{M} \left(\mathbf{e}_{\mathcal{A}_{t_2}} - \mathbf{l}_j \mathbf{s}_j^{(A_{t_2})} \right). \quad (10)$$

This formulation adjusts a particular column of \mathbf{L} and all non-zero corresponding entries of the $\mathbf{s}^{(t)}$'s. Additionally, instead of using squared-loss as our objective, we are free to choose \mathbf{M} and \mathbf{W} in a manner to weight certain errors, either between different features or between different tasks, more heavily. We can think of \mathbf{M} as analogous to $\mathbf{D}^{(t)}$ in Equation 7, except that \mathbf{M} must be shared among all tasks that share a particular latent model component and cannot be set independently for each task as in Equation 7. Additionally, the off-diagonal terms of the matrix \mathbf{W} correspond to extra terms that do not appear in Equation 7; these additional terms enforce relationships between tasks. Essentially, we can think of \mathbf{M} as a feature relationship matrix and \mathbf{W} as a task relationship matrix.

While there are many ways to choose \mathbf{M} and \mathbf{W} (we leave an exploration of these other options as future work), here we begin by setting all off-diagonal elements of \mathbf{W} to 0 in order to remove the inter-task penalty terms. We set \mathbf{M} as a consensus of each of the $\mathbf{D}^{(t)}$'s by computing their mean: $\mathbf{M} = \frac{1}{|\mathcal{A}|} \sum_{t \in \mathcal{A}} \mathbf{D}^{(t)}$. Next, we set the value of the diagonal entries of \mathbf{W} such that for each task we rescale \mathbf{M} to closely approximate the corresponding $\mathbf{D}^{(t)}$. Specifically, we set the t th diagonal element of \mathbf{W} as:

$$w_{t,t} = \frac{\mathbf{1}^\top \mathbf{D}^{(A_t)} \mathbf{1}}{\sum_{t' \in \mathcal{A}} \mathbf{1}^\top \mathbf{D}^{(t')} \mathbf{1}}, \quad (11)$$

where $\mathbf{1}$ is the vector of all 1's of the appropriate dimensionality (i.e., $\mathbf{1}^\top \mathbf{D}^{(t)} \mathbf{1}$ sums all entries of $\mathbf{D}^{(t)}$).

To be suitable for use with GSVD, both \mathbf{M} and \mathbf{W} must be PSD. \mathbf{M} is guaranteed to be PSD since it is an average of PSD matrices (the individual $\mathbf{D}^{(t)}$'s). \mathbf{W} is PSD since it is a diagonal matrix with all non-negative diagonal entries (the diagonal entries are non-negative since they are each a summation of the elements of a PSD matrix which is guaranteed to be positive).

Now that we have presented the modifications related to the dictionary update step of K-SVD, we write the updated method for computing $\mathbf{s}^{(t)}$ (originally defined

Algorithm 2 MTL-SVD

input training data $(\mathbf{X}^{(1)}, \mathbf{y}^{(1)}), \dots, (\mathbf{X}^{(T)}, \mathbf{y}^{(T)})$;
 dictionary size k
 init \mathbf{L} using random column vectors of unit length
for $i \in \{1, \dots, T\}$ **do**
 $(\boldsymbol{\theta}^{(i)}, \mathbf{D}^{(i)}) \leftarrow \text{singleTaskLearner}(\mathbf{X}^{(i)}, \mathbf{y}^{(i)})$
end for
loop until convergence **do**
 for $t \in \{1, \dots, T\}$, perform update in Eqn. 12
 for $j \in \{1, \dots, k\}$, perform update in Eqn. 9
end loop
return \mathbf{L}

in Equation 2) as:

$$\mathbf{s}^{(i)} \leftarrow \arg \min_{\mathbf{s}} \{ \|\mathbf{L}\mathbf{s} - \mathbf{x}_i\|_{\mathbf{D}^{(i)}}^2 + \mu \|\mathbf{s}\|_0 \}, \quad (12)$$

which we solve using the Lasso (Tibshirani, 1996). The complete MTL-SVD algorithm is given in Algorithm 2.

3.3. Computational Complexity

First, MTL-SVD uses a single task learner to compute the tuples $(\boldsymbol{\theta}^{(t)}, \mathbf{D}^{(t)})$ for all tasks. We use the function $\xi(d, n_t)$ to represent the computational complexity of the base learning algorithm on a problem of dimensionality d with n_t training instances. Next, the update of $\mathbf{s}^{(t)}$ given in Equation 12 can be solved using the LASSO algorithm in time $O(d^3 + d^2k + k^2d)$. Each update of a particular latent basis component and the corresponding entries in \mathbf{S} (i.e., Equations 9, 5, and 6) involves computing the square root of a $d \times d$ dimensional matrix as well as taking the SVD of a $d \times r$ dimensional matrix (where r is the maximum number of tasks that use a single latent basis vector). These two steps yield a complexity of $O(d^3 + r^2d)$. We must repeat the updates in Equations 9, 5, and 6 for each entry of $\mathbf{s}^{(t)}$. After an initial startup cost to compute the single task models of $O(\sum_{t=1}^T \xi(d, n_t))$, the per-iteration complexity of MTL-SVD is $O(d^2k + k^2d + kd^3 + kr^2d)$. In comparison, GO-MTL (Kumar & Daumé III, 2012), a current state-of-the-art multi-task learning algorithm, has a per iteration computational complexity of $O(d^3k^3)$.

4. Lifelong Learning Using K-SVD

Next, we show how to adapt the MTL-SVD algorithm to a lifelong learning setting in which new tasks arrive in an online fashion. We call our new algorithm *ELLA-SVD*, since it is an alternate formulation of ELLA.

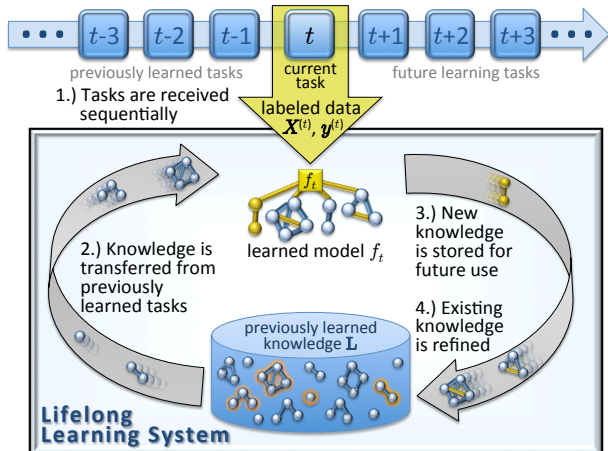


Figure 1. An illustration of the lifelong learning process.

4.1. The Lifelong Learning Problem

We now consider a variant of the multi-task learning setting in which the agent faces a *series* of supervised learning tasks $\mathcal{Z}^{(1)}, \mathcal{Z}^{(2)}, \dots, \mathcal{Z}^{(T_{\max})}$ (Figure 1). In contrast to the MTL setting, we assume that the learner receives the tasks in an online fashion and *a priori* does not know the total number of tasks T_{\max} , the distribution of these tasks, or their order.

Each time step, the agent receives a batch of labeled training data for some task t , either a new task or as additional training data for a previously learned task. In the lifelong learning setting, we use T to denote the number of tasks the agent has encountered so far, with $0 \leq T \leq T_{\max}$. After receiving each batch, the agent may be asked to make predictions on data instances of any previous task. Its goal is to construct task models $f^{(1)}, \dots, f^{(T)}$ where each $f^{(t)} : \mathbb{R}^d \mapsto \mathcal{Y}^{(t)}$ such that: (1) each $f^{(t)}$ will approximate $\hat{f}^{(t)}$ to enable the accurate prediction of labels for new instances, (2) each $f^{(t)}$ can be rapidly updated as the agent encounters additional training data for known tasks, and (3) new $f^{(t)}$'s can be added efficiently as the agent encounters new tasks. We assume that the total number of tasks T_{\max} and the total number of data instances $\sum_{t=1}^{T_{\max}} n_t$ will be large, and so a lifelong learning algorithm must have a computational complexity to update the task models that scales favorably with both quantities. Figure 1 illustrates the lifelong learning process.

4.2. Adaptions to MTL-SVD to Allow Efficient Lifelong Learning

The proposed MTL-SVD algorithm is inapplicable to the lifelong learning setting, since as each new task is presented to the learner, MTL-SVD would need to be rerun until convergence—a process could take arbitrarily long. As in earlier work (Ruvolo & Eaton, 2013),

we employ an update strategy that, where upon receiving data for task t , only the corresponding entry $\mathbf{s}^{(t)}$ is updated. Following this update, \mathbf{L} is then adapted to optimize performance across *all* tasks seen so far; however, in contrast to MTL-SVD these two steps are not repeated until convergence. Instead, they are each performed once for each new batch of training data. Despite these simplifications, we show that there is a small performance penalty incurred (both theoretically and empirically). This allows us to make the following modifications to the MTL-SVD algorithm to enable it to operate in a lifelong learning setting:

1. When receiving training data for task t , the update in Equation 12 (for updating the value of the $\mathbf{s}^{(t)}$'s) is only performed for task t .
2. The updates in Equations 9, 5, and 6 are performed only on the subset of columns of \mathbf{L} corresponding to the non-zero entries of $\mathbf{s}^{(t)}$ (i.e., we only update the latent basis vectors that are used to construct the model for the current task).
3. The two steps of the K-SVD algorithm are not repeated until convergence, but instead each performed only once per new batch of training data.

These modifications provide large gains in efficiency for three reasons. First, we do not need to update the $\mathbf{s}^{(t)}$'s for all previous tasks (which becomes expensive as T grows large). Second, the GSVD update step only has to be performed a number of times equal to the number of non-zero entries in $\mathbf{s}^{(t)}$ (which is a large savings when $\mathbf{s}^{(t)}$ is sparse). Third, we do not iterate updates of $\mathbf{s}^{(t)}$ and \mathbf{L} until convergence; instead we only need to run each step once.

4.3. Computational Complexity

As in Section 3.3, we use the function $\xi(d, n_t)$ to represent the computational complexity of the base learning algorithm on a problem of dimensionality d with n_t training instances. Next, the update of $\mathbf{s}^{(t)}$ given in Equation 12 can be solved using the LASSO algorithm in time $O(d^3 + d^2k + k^2d)$. Each update of a particular latent basis component and the corresponding entries in \mathbf{S} (i.e., Equations 9, 5, and 6) involves computing the square root of a $d \times d$ dimensional matrix as well as taking the SVD of a $d \times r$ dimensional matrix (where r is the number of tasks that utilize the latent basis component being updated). These two steps yield a complexity of $O(d^3 + r^2d)$. We must repeat the updates in Equations 9, 5, and 6 for each non-zero entry of $\mathbf{s}^{(t)}$. If we use q to indicate the number of such non-zero entries, then we arrive at a total complexity for incorporating a new batch of training data of $O(\xi(d, n_t) + d^2k + k^2d + qd^3 + qr^2d)$. In comparison, ELLA (Ruvolo & Eaton, 2013) has a computational

complexity for incorporating a new batch of training data of $O(\xi(d, n_t) + d^3k^2)$, which is significantly less efficient than our proposed ELLA-SVD.

5. Extensions

In this section, we modify the original ELLA (Ruvolo & Eaton, 2013) to include an incremental update, which reduces its complexity to more closely match ELLA-SVD. We also show that this incremental update can be integrated into ELLA-SVD to yield a hybrid approach. In the next section, we show that this hybrid approach has strong empirical performance.

5.1. ELLA with an Incremental Update

ELLA-SVD has a much lower computational complexity than the original ELLA. To facilitate better empirical comparison between the algorithms, we explore a modified version of ELLA, which we call *ELLA Incremental*, that closely matches the complexity of ELLA-SVD. We incorporate an incremental update into ELLA that updates each of the columns of \mathbf{L} independently when data for a new task is received. Further, only the columns of \mathbf{L} that are non-zero in the current $\mathbf{s}^{(t)}$ are updated and each column is updated at most once per batch of training data. The cost incurred for this more efficient update to \mathbf{L} is that ELLA Incremental is not guaranteed to achieve the globally optimal updated value of \mathbf{L} (given a fixed values of \mathbf{S}) as can be guaranteed in the original ELLA.

5.2. Hybrid Approach

Our proposed ELLA-SVD can be combined with ELLA Incremental into a hybrid approach. In this hybrid, we first perform one update step of ELLA-SVD and then one update step of ELLA Incremental. To guard against the two updates interfering with each other, if the ELLA-SVD step degrades the quality of our solution to Equation 8, then we only use the ELLA Incremental update for that iteration. We call this hybrid approach *ELLA Dual Update*. This approach maintains the strong advantage in computational complexity over ELLA of both ELLA-SVD and ELLA Incremental. Additionally, as we show in the next section, the hybrid approach performs much better than ELLA-SVD in situations where the assumptions of the ELLA-SVD algorithm are inaccurate.

6. Experiments

We evaluated four different algorithms for lifelong learning on four learning problems (one using synthetic data and the other three using real data). The four al-

gorithms that we evaluated were: (1) ELLA, as defined in (Ruvolo & Eaton, 2013), (2) ELLA-SVD, (3) ELLA Incremental, and (4) ELLA Dual Update.

We are primarily interested in developing a technique that closely approximates the accuracy of the ELLA algorithm, but with better computational complexity.

6.1. Data Sets

We tested each algorithm on four multi-task data sets: (1) synthetic regression tasks, (2) student exam score prediction, (3) land mine detection from radar images, and (4) facial expression recognition.

Synthetic Regression Tasks We created a set of $T_{\max} = 100$ random tasks with $d = 13$ features and $n_t = 100$ instances per task. The task parameter vectors $\boldsymbol{\theta}^{(t)}$ were generated as a linear combination of $k = 6$ randomly generated latent components in \mathbb{R}^{12} . The vectors $\mathbf{s}^{(t)}$ had a sparsity level of 0.5 (i.e., half the latent components were used to construct each $\boldsymbol{\theta}^{(t)}$). The training data $\mathbf{X}^{(t)}$ was generated from a standard normal distribution. The training labels for each task were given as $\mathbf{y}^{(t)} = \mathbf{X}^{(t)\top} \boldsymbol{\theta}^{(t)} + \epsilon$, where each element of ϵ is independent noise generated from a standard normal distribution. A bias term was added as the 13th feature prior to learning.

London School Data The London Schools data set consists of exam scores from 15,362 students in 139 schools. We treat the data from each school as a separate task. The goal is to predict the exam score of each student. We use the same feature encoding as used by Kumar & Daumé III (2012), where four school-specific and three student-specific categorical variables are encoded as a collection of binary features. We use the exam year and a bias term as additional features, giving each data instance $d = 27$ features.

Land Mine Detection In the land mine data set (Xue et al., 2007), the goal is to detect whether or not a land mine is present in an area based on radar images. The input features are automatically extracted from radar data and consist of four-moment based features, three correlation-based features, one energy-ratio feature, one spatial variance feature, and a bias term; see (Xue et al., 2007) for more details. The data set consists of a total of 14,820 data instances divided into 29 different geographical regions. We treat each geographical region as a different task.

Facial Expression Recognition This data set is from a recent facial expression recognition challenge (Valstar et al., 2011). Each task involves recognizing one of three facial action units (#5: upper lid raiser, #10: upper lip raiser, and #12: lip corner pull) from

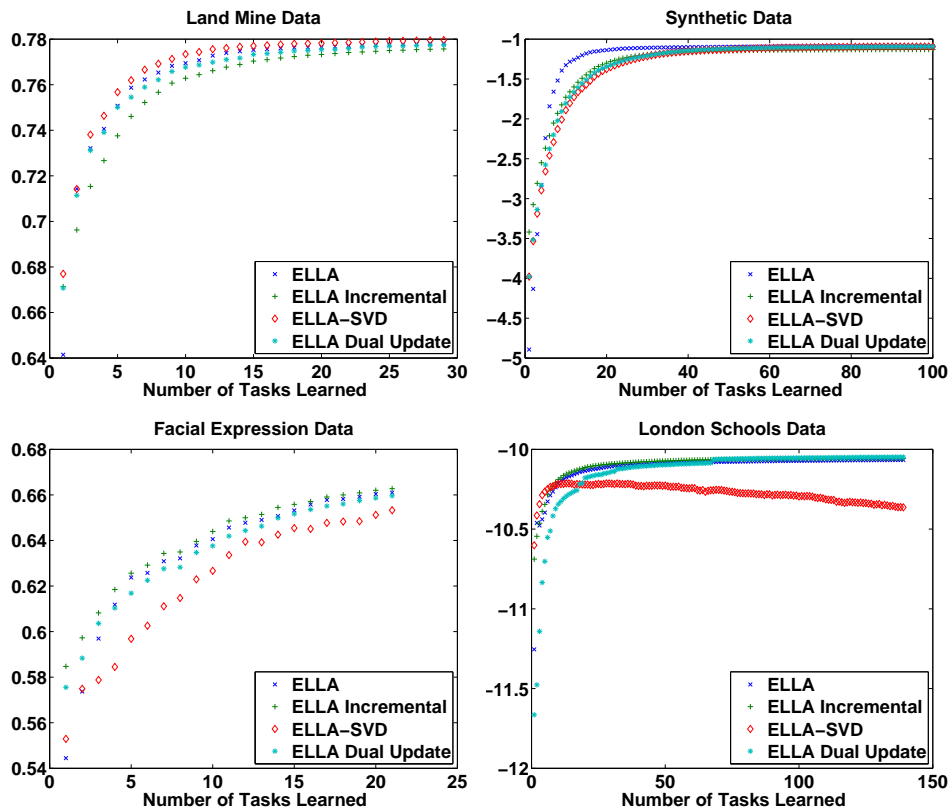


Figure 2. Learning curves for each of the four methods tested on four different datasets. Learning performance was evaluated after learning each task on *all* tasks.

an image of one of seven subjects’ faces. There are a total of 21 tasks, each with 450–999 images. We use the same feature encoding as Ruvolo & Eaton (2013), using a multi-scale Gabor pyramid to extract 2,880 Gabor features for each image, then reducing them to 100 dimensions using PCA, and adding a bias term.

6.2. Evaluation Procedure

Each task was presented sequentially to each algorithm according to the lifelong learning framework (Section 4.1) as a single batch of data that contained all training instances for that task. For each task, the training data was divided into both a training and a test set (with 50% of the data designated for each). The task order was also randomized. Each experiment was repeated 100 times to smooth out variability.

We generated learning curves for each method on each data set by averaging the performance across all tasks. We used the area under the ROC (AROC) as our metric for accuracy on classification tasks and negative root mean-squared (-rMSE) as our metric for regression tasks. We chose to use AROC rather than accuracy due to the skewed class distribution for the land

mine detection and facial expression recognition tasks. To evaluate the progress of learning performance as new tasks were presented, we measured the average task performance across all learned tasks. For tasks that were already learned this was straightforward: we simply evaluated how well the currently learned model for that task generalized to the corresponding test set. For tasks that had yet to be learned, we created a model for the task by fitting the currently learned basis \mathbf{L} to the training data for the task (but did not modify \mathbf{L}). Therefore, the learning curves evaluate both how well the current basis models previously learned tasks as well as how well it generalizes to *unseen* tasks.

The λ and k parameters were independently selected for each method to maximize the average (over all trials) final performance (i.e., the rightmost value on each learning curve). While fitting the parameters to maximize performance on the test set inflates the performance values compared to fitting these parameters on a validation set, the relative performance levels of the different algorithms (which is our principal focus) can still be compared. We conducted the grid-search over all combinations of $\lambda \in \{e^{-5}, \dots, e^5\}$ and $k \in \{1, \dots, 10\}$; μ was set to e^{-5} for all algorithms.

6.3. Results

The results of our evaluation are given in Figure 2. The proposed ELLA-SVD approach is better than all other methods on the land mine task. Specifically, the approach performs much better than the other efficient update approach, ELLA Incremental. On the synthetic regression tasks, the original ELLA method is clearly the best, with the ELLA-SVD and ELLA Incremental approaches lagging behind.

In contrast to the strong performance of ELLA-SVD on land mine and the synthetic tasks, ELLA-SVD does not perform well on either facial expression recognition or student exam score prediction. In particular, the performance of ELLA-SVD on student exam score prediction actually *declines* as it learns more tasks. Further investigation revealed that the cause of this problem was that the matrix \mathbf{M} formed as a consensus of the $\mathbf{D}^{(t)}$'s (which is required for Equation 9) is a poor approximation to the true objective function we would like to minimize (Equation 8). The primary reason for this poor approximation is that the input distributions for each task (i.e., each school) are quite different due to the school-specific features of each instance. In this case, the ELLA-SVD updates turn out to be counter-productive.

We proposed the ELLA Dual Update approach in order to get the best of all worlds. That is, we seek to achieve the high performance of ELLA-SVD on tasks where it is appropriate for application (e.g., for land mine detection), and to fallback to ELLA Incremental when ELLA-SVD performs poorly (e.g., for the London schools data). The results for the Dual Update version shown in Figure 2 suggest that this hybrid approach is successful. The performance of the ELLA Dual Update approach clusters tightly with the best performing algorithm for each learning problem (with the exception of the synthetic regression tasks, for which none of the more-efficient approaches does as well as the original ELLA).

7. Conclusion

We explored the use of the K-SVD algorithm in the lifelong machine learning setting. Adapting the approach of Aharon et al. (2006) to the lifelong learning setting required several key innovations including: (1) replacing the SVD step in the original algorithm with a generalized SVD, and (2) selectively updating components of the model as new task data is presented to the algorithm. We showed that this new algorithm, called ELLA-SVD, performs well on problems where the input distributions of the data are similar.

For domains where the input distributions are *not* similar, we showed that a hybrid approach (in which we interleave the ELLA-SVD update with another efficient update step called ELLA Incremental) performs robustly. In future work, we will conduct experiments to better understand the tradeoffs between ELLA-SVD and ELLA Incremental. Additionally, we plan to test our more-efficient versions of ELLA in settings where applying the original ELLA is computationally intractable (e.g., when k and d are large).

Acknowledgements

This research was supported by ONR grant N00014-11-1-0139. We thank Terran Lane, Diane Oyen, and the anonymous reviewers for their helpful feedback.

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