A Fully Tensorized Recurrent Neural Network

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Abstract
Recurrent neural networks (RNNs) are powerful tools for sequential modeling, but typically require significant overparameterization and regularization to achieve optimal performance. This leads to difficulties in the deployment of large RNNs in resource-limited settings, while also introducing complications in hyperparameter selection and training. To address these issues, we introduce a “fully tensorized” RNN architecture which jointly encodes the separate weight matrices within each recurrent cell using a lightweight tensor-train (TT) factorization. This approach represents a novel form of weight sharing which reduces model size by several orders of magnitude, while still maintaining similar or better performance compared to standard RNNs. Experiments on image classification and speaker verification tasks demonstrate further benefits for reducing inference times and stabilizing model training and hyperparameter selection.

1 Introduction
Recurrent neural networks (RNNs) represent a model family that is well-suited for tasks involving sequential data. Although early RNNs were limited by the problem of vanishing gradients during training, this was largely solved by the development of gated RNNs such as long short-term memory (LSTM) and gated recurrent unit (GRU) models (Hochreiter and Schmidhuber [1997], Cho et al. [2014]), which employ a collection of independent weight matrices to control the propagation of gradients. Such models have allowed RNNs to attain impressive performance in tasks such as speech recognition, language modeling, time series forecasting, and video classification.

RNNs typically employ large hidden states to achieve better performance in difficult modeling tasks, which in turn leads to a significant increase in parameters used to specify large weight matrices. The memory and compute issues associated with running such models, particularly in the limited setting of mobile and embedded devices, has led to the use of various techniques for model compression, including model distillation (Hinton, Vinyals, and Dean [2015]), alternate matrix decompositions (Sainath et al. [2013]), and quantization of network weights (He et al. [2019]). The use of such compression strategies is supported by the observation that standard representations of neural networks contain significant amounts of redundancy (Denil et al. [2013], Cheng et al. [2015]).

In this work we use the tensor-train (TT) formalism, a means of efficiently representing multi-modal tensors, to achieve significant compression of the model parameters associated with various RNN architectures. In contrast to previous work (Tjandra, Sakti, and Nakamura [2017], Yang, Krompass, and Tresp [2017]), we apply the TT formalism jointly to all weight matrices within the RNN, leading to a “fully tensorized” form of weight sharing, where various gate matrices are encoded within a single TT format. This permits the development of extremely lightweight, end-to-end trainable models, even in the presence of high-dimensional hidden states or input representations.

Experiments on image classification and speaker verification show that fully tensorized TT-RNNs give comparable or better performance relative to their uncompressed counterparts. We demonstrate that our method leads to state-of-the-art performance on the LibriSpeech dataset, producing a 16% reduction in speaker verification error while simultaneously allowing for a 200-fold compression in model parameters.

1.1 Notation
We use bold lower-case letters \( \mathbf{a} \) to denote vectors, bold upper-case letters \( \mathbf{W} \) to denote matrices, and bold calligraphic letters to denote tensors \( \mathbf{T} \). Tensor elements are indexed as \( \mathbf{a}(i) \), \( \mathbf{W}(i, j) \) and \( \mathbf{T}(i_1, i_2, ..., i_d) \), for the respective cases of vectors, matrices, and more general \( d \)’th-order tensors. The notation \( \mathbf{a} \ast \mathbf{b} \) represents the element-wise Hadamard product between vectors of equal size. The collection of integers \( \{1, 2, \ldots, n\} \) is denoted as \( [n] \).

2 Recurrent Neural Networks and the Tensor-Train Decomposition
In this section we first give an overview of common RNN architectures, before introducing the tensor-train decomposition and describing its use for “tensorizing” large weight matrices within neural networks.
2.1 RNN Architectures

Recurrent neural networks (RNN) define a paradigm for learning from sequential data. The recurrent unit of an RNN defines an iterative procedure whose outputs and hidden state at each time step \( t \) are a non-linear function of \( x_t \), the input at \( t \), and \( h_{t-1} \), the hidden state at time \( t-1 \). Many different functions have been proposed for this nonlinear recurrent unit, and we describe two representative choices, long short-term memory (LSTM) and gated recurrent unit (GRU).

Long Short-Term Memory

The LSTM cell uses three "gates" to control the flow of information, and divides its hidden state into a memory cell state \( c \) and regular hidden state \( h \), of identical dimension \( D \). These are jointly updated as

\[
\begin{align*}
    c_t &= u_t \ast \tilde{c}_t + f_t \ast c_{t-1} \\
    h_t &= o_t \ast \tanh(c_t),
\end{align*}
\]

where the candidate cell state (\( \tilde{c}_t \)), update gate (\( u_t \)), forget gate (\( f_t \)), and output gate (\( o_t \)) vectors are given by

\[
\begin{align*}
    \tilde{c}_t &= \tanh(W^{(c)} x_t + U^{(c)} h_{t-1} + b^{(c)}) \\
    u_t &= \sigma(W^{(u)} x_t + U^{(u)} h_{t-1} + b^{(u)}) \\
    f_t &= \sigma(W^{(f)} x_t + U^{(f)} h_{t-1} + b^{(f)}) \\
    o_t &= \sigma(W^{(o)} x_t + U^{(o)} h_{t-1} + b^{(o)}).
\end{align*}
\]

In the above, \( x_t \in \mathbb{R}^M \) and \( h_t \in \mathbb{R}^D \) are the input and hidden state vectors respectively, while \( W^{(c)}, W^{(u)}, W^{(f)}, W^{(o)} \in \mathbb{R}^{D \times M} \) are the input-hidden transition matrices, and \( U^{(c)}, U^{(u)}, U^{(f)}, U^{(o)} \in \mathbb{R}^{D \times D} \) are the hidden-hidden transition matrices.

Gated Recurrent Unit

The GRU is defined by two (update and relevance) gates and a single hidden state

\[
\begin{align*}
    h_t &= u_t \ast \tilde{h}_t + (1-u_t) \ast h_{t-1},
\end{align*}
\]

where

\[
\begin{align*}
    \tilde{h}_t &= \tanh(W^{(h)} x_t + U^{(h)} (r_t \ast h_{t-1}) + b^{(h)}) \\
    u_t &= \sigma(W^{(u)} x_t + U^{(u)} h_{t-1} + b^{(u)}) \\
    r_t &= \sigma(W^{(r)} x_t + U^{(r)} h_{t-1} + b^{(r)}).
\end{align*}
\]

The number of parameters for either of the above RNN units is \( gD(M + D) \), where \( g \) is the number of distinct gates, which is 4 for an LSTM and 3 for a GRU. Given any factorization of the input and hidden dimensions into positive integers as \( D = \prod_{k=1}^n d_k \) and \( M = \prod_{k=1}^n m_k \) (where \( d_k, m_k \geq 1 \)), this parameter count can be expressed as

\[
N_{\text{dense}} = gD(M + D) = g \left( \prod_{k=1}^n d_k m_k + \prod_{k=1}^n d_k \right) = O(d^n (m^n + d^n)),
\]

where \( d = \max_k d_k \) and \( m = \max_k m_k \). This version of the parameter count will allow for an easier comparison of typical RNN models with the tensorized RNNs introduced below.

2.2 Tensor-Train Decomposition

The tensor-train (TT) decomposition, introduced in [Oseledets 2011] and equivalent to the earlier matrix product state model of many-body physics [Vidal 2003], gives a method for representing higher-order tensors as a type of iterated low-rank factorization. A TT representation of an \( n \)-th order tensor \( T \in \mathbb{R}^{p_1 \times p_2 \times \cdots \times p_n} \) is a tuple of \( n \) tensors \( G_{[n]} = (G_1, G_2, \ldots, G_n) \), called the TT cores. Each core has dimension \( G_k \in \mathbb{R}^{p_k \times r_k \times r_k} \), where the \( r_k \) for \( k \in \{1, \ldots, n-1\} \) are hyperparameters called the TT ranks of the model. Given a collection of TT cores, the tensor \( T \) associated with these cores has elements given by the following vector-matrix-vector products

\[
T(i_1, i_2, \ldots, i_n) = G_1(i_1) G_2(i_2) \cdots G_d(i_n),
\]

where \( G_k(i_k) = G_k(i_k, \ldots) \in \mathbb{R}^{r_k \times r_k} \) indicates an index-dependent matrix associated with the \( k \)-th core, with each \( i_k \in [p_k] \) and \( r_0, r_n \) each taken to be 1. We will refer to \( T \) as the "global" tensor encoded by the TT cores, which constitute a "local" representation of \( T \).

The TT decomposition is capable of exactly representing any \( n \)-th order tensor given sufficiently large TT ranks using the TT-SVD procedure of [Oseledets 2011], but a more common practice is to fix the TT ranks at small values and use the core tensors as a compact parameterization which is optimized to minimize some loss function defined on the global tensor. This approach is not limited to cases where higher-order tensors are already present, as any vector \( v \) with dimension \( P = \prod_{k=1}^n p_k \) can be reshaped into an \( n \)-th order tensor \( V \in \mathbb{R}^{p_1 \times \cdots \times p_n} \). Such "TT vectors" provide an efficient description requiring only \( \sum_{k=1}^n p_k r_{k-1} r_k \) parameters when all TT ranks \( r_k \) and core dimensions \( p_k \) are bounded, compared with \( P \) parameters for a dense representation.

The same procedure can be applied to matrices of shape \( D \times M \) when \( D = \prod_{k=1}^n d_k \) and \( M = \prod_{k=1}^n m_k \), yielding a TT matrix defined by \( n \) tensor cores. In this case we choose each TT core \( G \) to have four indices with respective dimensions \( d_k, m_k, r_{k-1} \), and \( r_k \), and denote the associated index-dependent matrices by \( G_k(i_k, j_k) = G_k(i_k, j_k, \ldots) \in \mathbb{R}^{r_k \times r_k} \), for \( i_k \in [d_k] \) and \( j_k \in [m_k] \).

2.3 Tensorizing Neural Networks

The bulk of the parameters in a neural network consist of large weight matrices represented in dense format. It was shown in [Novikov et al. 2013] that the representation of these matrices as TT matrices allows for a significant reduction in parameter count, while introducing little or no additional error in the performance of the network.

Given a weight matrix \( W \) of shape \( D \times M \), where \( D = \prod_{k=1}^n d_k \) and \( M = \prod_{k=1}^n m_k \), then the affine transformation implemented as part of a typical neural network layer takes the form \( y = Wx + b \). In a tensorized neural network, \( x, y, b \) are represented normally as dense vectors, while the weight matrix \( W \) is represented in TT format. The affine transformation is carried out by first using multilinear tensor contractions to perform the multiplication \( Wx \), with \( x \) reshaped into a dense \( n \)-th order tensor \( X \), and then using standard
dense addition for the bias vector $b$. The output vector $y$ can be described in reshaped form as the tensor $\mathcal{Y}$ with elements

$$
\mathcal{Y}(i_1, \cdots, i_d) = \mathcal{B}(i_1, \cdots, i_d) + \sum_{j_1, \cdots, j_n} (\mathcal{G}_1(i_1, j_1) \cdots \mathcal{G}_d(i_d, j_d)) \mathcal{X}(j_1, \cdots, j_d).
$$

(7)

By carrying out the above summations (including those implicit in the matrix-vector products) in an optimal order, (7) can be evaluated with a total cost of $O(nr^2dM)$, where $r = \max_k r_k$. In the typical setting where $r$, $m$, and $d$ remain bounded as $D$ and $M$ are increased, this cost is $O(\log(\max(D, M))M)$, compared to $O(DM)$ for the usual affine map. This representation is also compact, requiring only $O(nr^2dM) = O(\log(\max(D, M)))$ parameters, compared to $O(DM)$ parameters for a dense representation.

For clarity, we refer to a fully-connected layer represented in tensor-train form as a tensor-train layer (TTL), and denote the linear portion of the operation implemented in (7) as TTL($x; \mathcal{G}_{[n]}$).

### 2.4 Tensorizing RNNs

We describe a straightforward application of the above tensorization procedure to LSTM models, as utilized in [Takeda, Sakti, and Nakamura 2017, Yang, Krompass, and Tresp 2017], which allows for a significant reduction in the model parameter count. In the next section we propose an extension of this procedure which permits an even greater degree of compression to be attained.

An LSTM recurrent unit contains 8 weight matrices, each providing contributions to one of the four independent gate vectors coming from an input vector $x_t$ or previous hidden vector $h_t$. When these matrices are replaced by tensor-train matrices, (3) can be re-written as

$$
\begin{align*}
\tilde{c}_t &= \tanh(\text{TTL}(x_t; \mathcal{G}_{[n]}^{Wc}) + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{Uc}) + b^{(c)}) \\
u_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{Wu}) + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{Uu}) + b^{(u)}) \\
f_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{Wf}) + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{Uf}) + b^{(f)}) \\
o_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{Wo}) + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{Uo}) + b^{(o)}).
\end{align*}
$$

(8)

Each of the 8 weight matrices $V_e$ (where $V$ is one of $W$, $U$, and $e$ is one of $c$, $u$, $f$, or $o$) is replaced by its own collection of tensor-train cores $\mathcal{G}_{[n]}^{Ve}$, and we assume for simplicity that the same factorization of $D = \prod_{k=1}^n d_k$ and $M = \prod_{k=1}^n m_k$ is used for each of the 8 tensor-train matrices.

For a tensorized gated RNN with $g$ gates and an identical factorization for each tensor-train matrix, such as the LSTM above, the total parameter count is

$$
N_{\text{TTL}} = g \sum_{k=1}^n r_{k-1}r_kd_k(m_k + d_k) = O(gnr^2d(m + d)).
$$

(9)

Although the exact comparison of this count to (5) depends on the TT ranks $r_k$ and the number of cores $n$ employed, it is clear that for the typical case where $r, d, m \ll \min(M, D)$, a tensorized RNN will require significantly fewer parameters. However, the use of a separate TT matrix for each gate in the RNN unit still leads to a multiplicative factor of $g$ in (9).

### 3 Fully Tensorized RNNs

We now introduce a different tensorization method, where a tensor-train factorization is applied to entire collections of concatenated weight matrices, rather than to individual matrices. The efficient nature of the tensor-train decomposition leads to a further reduction in model parameters, with LSTMs requiring approximately four times fewer parameters compared to the tensorization above. We show more generally that gated RNNs with $g$ gates exhibit a roughly $g$-fold reduction in the parameter count with this method, on top of the already sizable reduction coming from the use of tensor-train matrices.

#### 3.1 Gate Concatenation

We achieve further compression of our tensorized RNN by jointly tensorizing the input-hidden weights, as well as the hidden-hidden weights. Taking the LSTM as an example, we first take the row-wise concatenation of the four input-hidden matrices $W^{(c)}$, $W^{(u)}$, $W^{(f)}$, $W^{(o)} \in \mathbb{R}^{D \times M}$, which gives a single input-hidden matrix $W \in \mathbb{R}^{4D \times M}$. More concretely, the concatenated weight matrices utilized are

$$
W = [W^{(c)}, W^{(u)}, W^{(f)}, W^{(o)}]^T, \\
U = [U^{(c)}, U^{(u)}, U^{(f)}, U^{(o)}]^T.
$$

(10)

For regular LSTMs with dense weight matrices, this concatenation gives a means of replacing four separate matrix-vector multiplications by a single larger multiplication, permitting greater parallelism. After the single vector $Wx$ is computed it can be split into four equal-sized pieces, each holding the value of one of the gate vectors.

When the concatenated weight matrices are represented as a tensor-train layer, this leads to the revised LSTM gate equations,

$$
\begin{align*}
\tilde{c}_t &= \tanh(\text{TTL}(x_t; \mathcal{G}_{[n]}^{W})_1 + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{U})_1 + b^{(c)}) \\
u_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{W})_2 + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{U})_2 + b^{(u)}) \\
f_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{W})_3 + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{U})_3 + b^{(f)}) \\
o_t &= \sigma(\text{TTL}(x_t; \mathcal{G}_{[n]}^{W})_4 + \text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{U})_4 + b^{(o)}),
\end{align*}
$$

(11)

where $\text{TTL}(x_t; \mathcal{G}_{[n]}^{W})_i$ and $\text{TTL}(h_{t-1}; \mathcal{G}_{[n]}^{U})_i$ are the $i$th equally-sized vectors in the TT matrix-vector products associated with $W$ and $U$, which contribute to the $c$, $u$, $f$, and $o$ gates. This process can be carried out analogously for a gated RNN with $g$ gates, where the matrices $W \in \mathbb{R}^{gD \times M}$, $U \in \mathbb{R}^{gD \times L}$ are each concatenations of $g$ separate matrices. An example of this process for a GRU model is given in Figure 4.
When tensorizing the individual weight matrices of an RNN in Section 2.4, the hidden and input dimensions were factored into smaller terms, as $D = \prod_{k=1}^{n} d_k$ and $M = \prod_{k=1}^{n} m_k$. For the case of concatenated weight matrices $W$ and $U$, a closely related factorization can be employed, namely $gD = \prod_{k=0}^{n} d_k$ and $M = \prod_{k=0}^{n} m_k$, where we take $d_0 = g$ and $m_0 = 1$, along with identical $d_k, m_k$ for all $k \geq 1$.

Taking $W$ as an example, a tensor-train decomposition relative to this augmented factorization will give the collection of $n + 1$ cores $G_{[n+1]}^{(W)} = (G_0^{(W)}, G_1^{(W)}, \ldots, G_n^{(W)})$, where the cores $G_k^{(W)}$ for $k > 1$ are shaped identically to a tensor-train factorization of any one of the single-gate weight matrices. The single new core appearing in this decomposition has a shape of $G_0^{(W)} \in \mathbb{R}^{g \times 1 \times 1 \times r_0}$, for a new TT rank parameter $r_0$, and removing the singleton indices gives a matrix $V^{(W)} \in \mathbb{R}^{g \times r_0}$. This leads to a revised parameter count

$$N_{TT} = gr_0 + \sum_{k=1}^{n} r_k d_k (m_k + d_k) = \mathcal{O}(n r^2 d (m + d)),$$

(12)

giving a compression ratio approximately $g$ times greater than (9). Using an example model in Table 1, we illustrate the level of compression and speedup in inference time that can be obtained for different configurations of our fully tensorized RNNs. This shows particular promise for the application of RNN models in settings with limited resources, such as edge devices. Finally, the training time for TT-RNNs is comparable to untensorized RNNs, although with a clear dependence on the TT rank.

3.3 Weight Sharing

Some intuition for this parameter reduction can be gained by interpreting the concatenated global matrix $W$ encoded by the TT cores $G_{[n+1]}^{(W)}$ in terms of the small matrix $V^{(W)}$ coming from the first core $G_0^{(W)}$. Seen this way, the contraction of the remaining TT cores $(G_1^{(W)}, \ldots, G_n^{(W)})$ gives a tensor which encodes a family of $r_0$ matrices $M_\alpha \in \mathbb{R}^{D \times M}$, jointly represented in TT format, while the top core $G_0$ acts as a matrix assigning each of the GRU gate matrices to a linear mixture of the TT matrices $M_\alpha$ (see Equation (3.3)).

4 Experiments

We benchmark the performance of TT-RNN models using experiments on image classification and speaker verification tasks. Results for TT-LSTM are reported here, while those for TT-GRU can be found in the supplementary material. Beyond assessing the accuracy in these tasks, we characterize trade-offs between compression and accuracy arising from different choices of TT rank and core layout. In the process, we find that the tensor-train parameterization acts as a form of regularization, leading to improved stability and generalization during training.

For simplicity and ease of comparison, all models in the following are trained without explicit regularization such as dropout, weight decay, or gradient clipping. The tensorized models were written in PyTorch (Paszke et al. 2019) using the tensor-train implementation from (Khrulkov et al. 2019), and are available on GitHub.\(^3\)

\(^3\)https://github.com/onucharles/tensorized-rnn
Table 1: Comparison of model size and per-step training and inference times of RNNs and TT-RNNs. Each model has a single recurrent layer with hidden size of 512, a linear projection layer of embedding size 256, and input dimension of 4,096. Each TT-RNNs has 2 cores, and \( r \) denotes the TT rank. For both LSTM and GRU models, the tensorized versions achieve significant compression of model parameters while reducing the inference time and, for smaller values of \( r \), decreasing the training time. All reported times were obtained on an Intel(R) Xeon(R) CPU E5-1650 v3 @ 3.50GHz with 128GB of RAM, and averaged over 100 runs.

| Model  | \( r \) | # params | Train time (s) | Eval. time (s) |
|--------|--------|----------|----------------|----------------|
| LSTM   | –      | 9,570,560| 12.84 ± 1.17   | 3.70 ± .19     |
| TT-LSTM| 2      | 21,248   | 9.37 ± .11     | 2.13 ± .13     |
|        | 3      | 30,720   | 11.92 ± .22    | 2.23 ± .13     |
|        | 4      | 40,192   | 15.55 ± .37    | 2.48 ± .25     |
| GRU    | –      | 7,212,288| 10.12 ± .26    | 2.53 ± .07     |
| TT-GRU | 2      | 19,200   | 8.09 ± .21     | 1.43 ± .09     |
|        | 3      | 27,136   | 9.18 ± .15     | 1.59 ± .08     |
|        | 4      | 35,072   | 11.23 ± .30    | 1.80 ± .10     |

### 4.1 Permuted Pixel MNIST

We first evaluate the TT-LSTMs on the permuted sequential MNIST task (LeCun, Cortes, and Burges 1998) in which the \( 28 \times 28 \) pixel images of handwritten digits are randomly rearranged using a fixed permutation into sequences of length 784. These are split into 50k training, 10k validation, and 10k test images, with the validation dataset used to determine the end of training by early stopping.

The LSTM and TT-LSTM were each chosen as single-layer models with 256 hidden units. Training was performed with a batch size of 256 and Adam optimizer, using a piecewise constant learning rate starting at 0.001.

Table 2 reports the digit classification accuracy, where the hidden dimensions of the TT-LSTMs are factored into either 2 or 3 TT cores using TT ranks of 2, 4, or 6 to connect adjacent cores. Although a clear tradeoff is present between compression and accuracy, even the largest TT-LSTM utilizes 46 times fewer parameter in total, while achieving comparable performance to the LSTM baseline (−0.28% classification accuracy).

Table 2: Comparison of TT-RNN and standard RNN models on the permuted pixel MNIST task. The models use a single-layer containing \( D = 256 \) hidden units, and are trained identically. The performance of the TT-RNNs varies with the parameter count, but achieves comparable accuracy to a standard RNNs while maintaining a compression ratio of 46 times and 25 times fewer parameters, in the TT-LSTM and TT-GRU respectively.

| Model     | Cores | \( r \) | #Params | Compr. | Acc. (%) |
|-----------|-------|--------|---------|--------|----------|
| LSTM      | –     | –      | 266,762 | –      | 89.77    |
| TT-LSTM   | 2     | 4      | 3,434   | 78     | 87.98    |
|           | 4     | 5      | 5,834   | 46     | 89.49    |
|           | 6     | 8      | 8,234   | 32     | 89.22    |
| GRU       | –     | –      | 201,482 | –      | 91.49    |
| TT-GRU    | 2     | 4      | 3,574   | 25     | 87.99    |
|           | 4     | 5      | 5,802   | 35     | 89.29    |
|           | 6     | 7      | 7,930   | 25     | 90.26    |
|           | 2     | 4      | 3,434   | 54     | 88.90    |
|           | 6     | 5      | 5,866   | 34     | 89.80    |

### 4.2 Speaker Verification

In the speaker verification problem, the objective is to ascertain if an utterance of speech belongs to a given individual, based on a collection of utterances labeled by individuals. We use the LibriSpeech dataset, containing around 1,000 hours of English language audiobook recordings (Panayotov et al. 2015), where training, validation, and testing are carried out on the train-clean-100, dev-clean, and test-clean partitions.

The TT-LSTM exhibits impressive performance in this more complex task, substantially improving on the LSTM baseline while using 200 times fewer parameters. The tensor-train parameterization of the TT-LSTM appears to be a form of implicit regularization, which leads to less overfitting while also minimizing the issue of vanishing and exploding gradients during training.

**Setup** Our model for speaker verification contains two main components, an utterance encoder and a similarity function, as in (Heigold et al. 2016; Xie et al. 2019). The utterance encoder consists of an RNN which computes fixed-dimensional embeddings from spectograms of input utterances, while the similarity function assigns similarity scores to pairs of embeddings.

We use the generalized end-to-end (GE2E) loss function (Wan et al. 2018) to train the model, which encourages embeddings of utterances to cluster based on the associated speaker. Given an embedding vector \( e_{ji} \) for the \( j \)th utterance by the \( j \)th speaker, the GE2E loss is

\[
L(e_{ji}) = -S_{ji,j} + \log \sum_{k=1}^{N} \exp(S_{ji,k}),
\]

where \( S_{ji,j} = w \cdot \cos(e_{ji}, e_k) + b \) is the scaled cosine similarity between the embedding \( e_{ji} \) and the centroid of the embeddings of speaker \( j \), denoted \( e_j \). The scaling coefficients \( w \) and \( b \) are initialized to 10 and −5 respectively. The full loss is then the sum of all utterance-specific losses, \( L = \sum_{j,i} L(e_{ji}) \).

We report performance in the speaker verification task using the equal error rate (EER) metric, which is the error rate on the receiver-operating characteristic (ROC) curve when the false positive rate and false negative rates are equal.

**Performance** Our utterance encoder consists of a single-layer LSTM with hidden size of 768, whose output is converted to an embedding of dimension 256 using a fully-connected linear layer. The input to this encoder is 40-bin \( \times \) 160-frame Mel spectograms of utterances. We compare reg-
Table 3: Performance of RNNs and TT-RNNs on the task of speaker verification. Models have a single layer with 768 hidden units and a linear projection layer of 256. The lowest ranked TT-RNNs outperform the RNNs on this more challenging task of speaker verification, achieving larger compression ratios of 653 (TT-LSTM) and 369 (TT-GRU). EER is the equal error rate (lower is better).

| Model   | Cores | r   | #Params  | Compr. | EER (%) |
|---------|-------|-----|---------|--------|---------|
| LSTM    | −     | −   | 2,682,114 | −      | 7.33    |
| TT-LSTM | 2     | 1   | 8,178   | 328    | 4.71    |
|         | 2     | 13,026 | 206    | 4.34   |
|         | 4     | 22,722 | 118    | 6.21   |
|         | 1     | 4,106  | 653    | 6.09   |
|         | 2     | 5,394  | 497    | 5.31   |
|         | 4     | 9,506  | 282    | 5.38   |
| GRU     | −     | −   | 2,063,106 | −      | 7.87    |
| TT-GRU  | 2     | 1   | 9,074   | 227    | 5.31    |
|         | 2     | 13,282 | 155    | 6.72   |
|         | 4     | 21,698 | 95     | 5.36   |
|         | 1     | 5,594  | 369    | 6.46   |
|         | 2     | 6,738  | 306    | 6.39   |
|         | 4     | 10,274 | 201    | 4.48   |

Using a standard LSTM in the encoder gives an EER of 7.33%, similar to the performance found in (Zhou et al. 2019). By contrast a TT-LSTM encoder led to significantly better EERs, with the best configuration achieving an EER of 4.34%. This increased accuracy was accompanied by a reduction in the total parameter count, from 2.6M parameters to only 13K. By reducing the TT rank, this parameter count can be further reduced while still maintaining higher accuracy than the LSTM baseline.

Analyzing the embeddings learned by the TT-LSTM further demonstrates the performance of the model in speaker verification. We use uniform manifold approximation and projection (UMAP) (McInnes, Healy, and Melville 2018) to project the 256-dimension embedding vectors into 2D space (Figure 2), which shows that the embeddings learned by the TT-LSTM effectively cluster the utterances from each speaker.

Regularization TT-LSTMs utilize a more compact set of weight parameters, which can be expressed as a low-dimensional family of weight matrices. To assess if this low-dimensional parameterization has benefits for regularization, we first examine the learning curves of TT-LSTMs and standard LSTMs during training (Figure 3 left). We observe that while LSTM encoders achieve lower loss during training, this loss is not reflected in the validation loss, likely due to overfitting. By contrast, the TT-LSTM shows better generalization, giving a smaller discrepancy between training and validation loss, and ultimately a lower validation EER.

To further test this generalization, we conduct the speaker verification experiments in a more data-limited setting, using between 20% and 100% of the training data. TT-LSTMs consistently performed better than the LSTM baseline when trained with small amounts of data (Figure 3 right).

Training Stability We observed during the initial hyperparameter search an increased robustness in the performance of TT-LSTMs relative to changes in the learning rate. Both LSTMs and TT-LSTM models were trained at a learning rate of 0.001, but increasing this to 0.01 led to an instability in the former and no noticeable impact on the latter. The distribution of gradients for this case is given in Figure 4. The standard LSTM exhibits vanishing gradients, effectively saturating at 0, while the gradients for TT-LSTM are distributed over a reasonable range.

5 Related Work

The compression of deep neural networks (DNNs) has been of interest for a long time. It has been shown that DNNs are typically parameterized in a redundant fashion, allowing the prediction of values of some parameters of a trained model given knowledge of the others (Denil et al. 2013).

Several approaches depend on some kind of post-processing after a large model has been trained. Model distillation (Ba and Caruana 2014; Hinton, Vinyals, and Dean 2015) for example is a successful technique which retrains a smaller model by using the output activations of the trained large model as labels, instead of the actual data labels. This was found to result in smaller models that are fast to train and match the performance of the larger models from which they were distilled. Quantization is another post-processing technique which uses a more coarse-grained representation for each parameter value, thereby reducing the memory needed to store a trained model’s parameters. One complication with these post-processing methods is that they are not end-to-end; the process of pruning the DNN is separate from training.

Matrix and tensor factorization techniques provide an alternative that is end-to-end trainable. A natural first step is...
to decompose parameter weight matrices in a low-rank matrix factorized format. This was done in (Sainath et al. 2013) to compress the last fully-connected layer of a convolutional neural network (CNN). Restricting to the last layer is limited in the compression achieved, since the other layers of the network themselves contain many parameters. However, utilizing this approach in internal layers results in a lower effective number of hidden units, ultimately hurting accuracy.

Tensor factorization methods, such as that employed here, can generally be used to decompose matrices in higher-dimensional space, and were used in (Yu et al. 2017) to capture higher order interactions in dynamical processes. The idea of tensorizing neural networks in an end-to-end trainable manner using tensor-train decomposition was first introduced in (Novikov et al. 2015), where a fully-connected layer was reshaped and factorized as a tensor in TT format to achieve impressive compression. The extension to convolutional layers was later given in (Garipov et al. 2016).

Different aspects of these ideas were extended to recurrent neural networks (RNN) in (Yang, Krompass, and Tresp 2017; Tjandra, Sakti, and Nakamura 2017). The work of (Yang, Krompass, and Tresp 2017) applied tensor-train layers to the large encoding matrices used for high-dimensional video input, allowing for simultaneous compression and improved performance in video classification. This was later followed by (Yin et al. 2020), which reported further gains through the use of the more complex Hierarchical Tucker decomposition in place of tensor trains.

By contrast, (Tjandra, Sakti, and Nakamura 2017) tensorized RNNs by assigning a separate TT matrix to each of the separate weight matrices in a recurrent cell, with a focus on GRU models. This allowed significant compression to be achieved not only with high-dimensional inputs, but also with high-dimensional hidden states. Our work is similar to (Tjandra, Sakti, and Nakamura 2017), but achieves further compression by jointly tensorizing the weights within each RNN cell. We show how this process leads to a novel form of weight sharing, which is verified experimentally to have tangible benefits for performance and compression. The use of a tensor-train parameterization is shown to represent an implicit regularization capable of improving training and generalization. Our TT-RNN model is available as open-source code, and can be used as a drop-in replacement for standard RNN models.

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