DeBERTa: Decoding-enhanced BERT with Disentangled Attention

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Abstract

Recent progress in pre-trained neural language models has significantly improved the performance of many natural language processing (NLP) tasks. In this paper we propose a new model architecture **DeBERTa** (**D**ecoding-**e**nhanced **BERT** with disentangled **a**ttention) that improves the BERT and RoBERTa models using two novel techniques. The first is the disentangled attention mechanism, where each word is represented using two vectors that encode its content and position, respectively, and the attention weights among words are computed using disentangled matrices on their contents and relative positions. Second, an enhanced mask decoder is used to replace the output softmax layer to predict the masked tokens for model pre-training. We show that these two techniques significantly improve the efficiency of model pre-training and performance of downstream tasks. Compared to RoBERTa-Large, a DeBERTa model trained on half of the training data performs consistently better on a wide range of NLP tasks, achieving improvements on MNLI by +0.9% (90.2% vs. 91.1%), on SQuAD v2.0 by +2.3% (88.4% vs. 90.7%) and RACE by +3.6% (83.2% vs. 86.8%). The DeBERTa code and pre-trained models will be made publicly available `https://github.com/microsoft/DeBERTa`.

1 Introduction

The Transformer has become the most effective neural network architecture for neural language modeling. Unlike recurrent neural networks (RNNs) that process text in sequence, Transformers apply self-attention to compute in parallel every word from the input text an attention weight that gauges the influence each word has on another, thus allowing for much more parallelization than RNNs for large-scale model training [1]. Since 2018, we have seen the rise of a set of large-scale Transformer-based Pre-trained Language Models (PLMs), such as GPT [2, 3], BERT [4], RoBERTa [5], XLNet [6], UniLM [7], ELECTRA [8], T5 [9], ALUM [10], StructBERT [11] and ERINE [12] . These PLMs have been fine-tuned using task-specific labels and created new state-of-the-art in many downstream natural language processing (NLP) tasks [13, 14, 15, 16, 17, 18].

In this paper, we propose a new Transformer-based neural language model **DeBERTa** (**D**ecoding-**e**nhanced **BERT** with disentangled **a**ttention) which has been proven to be more effective than RoBERTa and BERT as a PLM and after fine-tuning leads to better results on a wide range of NLP tasks.

DeBERTa makes two modifications to the BERT model. The first is the use of a disentangled attention mechanism for self-attention. Unlike BERT where each word in the input layer is represented using a vector which is the sum of its word (content) embedding and position embedding, each word in DeBERTa is represented using two vectors that encode its content and position, respectively, and the attention weights among words are computed using disentangled matrices based on their contents and relative positions, respectively. This is motivated by the observation that the attention weight of a

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word pair depends on not only their contents but their relative positions. For example, the dependency between the words "deep" and "learning" is much stronger when they occur next to each other than when they occur in different sentences.

Second, we enhance the output layer of BERT for pre-training. The output softmax layer of BERT is replaced with an Enhanced Mask Decoder (EMD) to predict the masked tokens during model pre-training. This is motivated by mitigating a mismatch between pre-training and fine-tuning. While fine-tuning, we use a task-specific decoder that takes the BERT output as input and produces the task labels. However, while pre-training we do not use any task-specific decoder and instead simply normalize the BERT output (logits) via softmax. We thus treat the pre-training task, masked language model (MLM), the same as any fine-tuning task, and add a task-specific decoder, which is implemented as a two-layer Transformer decoder and a softmax output layer, for pre-training.

We will show through a comprehensive empirical study that these two techniques substantially improve the efficiency of pre-training and the performance of downstream tasks. Compared to RoBERTa-Large, a DeBERTa model trained on half the training data performs consistently better on a wide range of NLP tasks, achieving improvements on MNLI by +0.9% (90.2% vs. 91.1%), on SQuAD v2.0 by +2.3%(88.4% vs. 90.7%), and RACE by +3.6% (83.2% vs. 86.8%).

2 Background

2.1 Transformer structure

A Transformer-based language model is composed of stacked Transformer blocks [1]. Each block contains a multi-head self-attention layer followed by a fully connected positional feed-forward network. The standard self-attention mechanism lacks a natural way to encode word position information. Thus, existing approaches add a positional bias to each input word embedding so that each input word is represented by a vector whose value depends on its content and position. The positional bias can be implemented using absolute position embedding [1, 2, 4] or relative position embedding [6, 19]. It has been shown that relative position representations are more effective for natural language understanding and generation tasks [20, 21]. The proposed Disentangled Attention mechanism differs from all existing approaches in that we represent each input word using two separate vectors that encode a word's content and position respectively, and attention weights among words are computed using disentangled matrices on their contents and relative positions.

2.2 Masked Language Model

Large-scale Transformer-based PLMs [4, 5, 22] are typically pre-trained on large amounts of text to learn contextual word representations using a self-supervision objective, known as Masked Language Model (MLM). Specifically, given a sequence $\boldsymbol{X} = \{x_i\}$, we corrupt it into $\tilde{\boldsymbol{X}}$ by masking 15% of its tokens at random and then train a language model parameterized by θ to reconstruct $\boldsymbol{X}$ by predicting the masked tokens $\tilde{x}$ conditioned on $\tilde{\boldsymbol{X}}$:

$$\max_{\theta} \log p_\theta(\boldsymbol{X}|\tilde{\boldsymbol{X}}) = \sum_{i \in \mathcal{C}} \log p_\theta(\tilde{x}_i = x_i|\tilde{\boldsymbol{X}}) \tag{1}$$

where $\mathcal{C}$ is the index set of the masked tokens in the sequence. The authors of BERT propose to keep 10% of the masked tokens unchanged, another 10% replaced with randomly picked tokens and the rest replaced with the [MASK] token. In DeBERTa, we use a MLM decoder for masked token prediction and replace the 10% masked but unchanged tokens with their position embedding vectors in the decoder.

3 Approach

3.1 Disentangled Attention

For a token at position i in a sequence, we represent it using two vectors, $\{\boldsymbol{H}_i\}$ and $\{\boldsymbol{P}_{i|j}\}$, which represent its content and relative position with the token at position j, respectively. The calculation of

the cross attention score between tokens i and j can be decomposed into four components as

$$A_{i,j} = \{\boldsymbol{H_i}, \boldsymbol{P_{i|j}}\} \times \{\boldsymbol{H_j}, \boldsymbol{P_{j|i}}\}^\mathsf{T}$$
$$= \boldsymbol{H_i}\boldsymbol{H_j}^\mathsf{T} + \boldsymbol{H_i}\boldsymbol{P_{j|i}}^\mathsf{T} + \boldsymbol{P_{i|j}}\boldsymbol{H_j}^\mathsf{T} + \boldsymbol{P_{i|j}}\boldsymbol{P_{j|i}}^\mathsf{T} \tag{2}$$

That is, the attention weight of a word pair can be computed as a sum of four attention scores using disentangled matrices on their contents and positions as *content-to-content*, *content-to-position*, *position-to-content*, and *position-to-position* [1].

Existing approaches [19, 21] to relative position encoding use a separate embedding matrix to compute the relative position bias in computing attention weights. This is equivalent to computing the attention weights using only the content-to-content and content-to-position terms in (2). We argue that the position-to-content term is also important since the attention weight of a word pair depends not only on their contents but on their relative positions, which can only be fully modeled using both the content-to-position and position-to-content terms. Since we use *relative* position embedding, the position-to-position term does not provide much additional information and is removed from (2) in our implementation.

Taking single-head attention as an example, the standard self-attention [1] can be formulated as:

$$\boldsymbol{Q} = \boldsymbol{H}\boldsymbol{W_q}, \boldsymbol{K} = \boldsymbol{H}\boldsymbol{W_k}, \boldsymbol{V} = \boldsymbol{H}\boldsymbol{W_v}, \boldsymbol{A} = \frac{\boldsymbol{Q}\boldsymbol{K}^\mathsf{T}}{\sqrt{d}}$$

$$\boldsymbol{H_o} = \mathrm{softmax}(\boldsymbol{A})\boldsymbol{V}$$

where $\boldsymbol{H} \in R^{N \times d}$ represents the input hidden vectors, $\boldsymbol{H_o} \in R^{N \times d}$ the output of self-attention, $\boldsymbol{W_q}, \boldsymbol{W_k}, \boldsymbol{W_v} \in R^{d \times d}$ the projection matrices, $\boldsymbol{A} \in R^{N \times N}$ the attention matrix, N the length of input sequence, and d the dimension of hidden state.

Denote k as the maximum relative distance, $\delta(i, j) \in [0, 2k)$ as the relative distance from token i to token j, which is defined as:

$$\delta(i,j) = \left\{ \begin{array}{rcl} 0 & \text{for} & i - j \leqslant -k \\ 2k - 1 & \text{for} & i - j \geqslant k \\ i - j + k & \text{others} & \end{array} \right. \tag{3}$$

We can represent the disentangled self-attention with relative position bias as (4), where $\boldsymbol{Q_c}, \boldsymbol{K_c}$ and $\boldsymbol{V_c}$ are the projected content vectors generated using projection matrices $\boldsymbol{W_{q,c}}, \boldsymbol{W_{k,c}}, \boldsymbol{W_{v,c}} \in R^{d \times d}$ respectively, $\boldsymbol{P} \in R^{2k \times d}$ represents the relative position embedding vectors shared across all layers (i.e., staying fixed during forward propagation), and $\boldsymbol{Q_r}$ and $\boldsymbol{K_r}$ are projected relative position vectors generated using projection matrices $\boldsymbol{W_{q,r}}, \boldsymbol{W_{k,r}} \in R^{d \times d}$, respectively.

$$\boldsymbol{Q_c} = \boldsymbol{H}\boldsymbol{W_{q,c}}, \boldsymbol{K_c} = \boldsymbol{H}\boldsymbol{W_{k,c}}, \boldsymbol{V_c} = \boldsymbol{H}\boldsymbol{W_{v,c}}, \boldsymbol{Q_r} = \boldsymbol{P}\boldsymbol{W_{q,r}}, \boldsymbol{K_r} = \boldsymbol{P}\boldsymbol{W_{k,r}}$$

$$\tilde{A}_{i,j} = \underbrace{\boldsymbol{Q_i^c}\boldsymbol{K_j^c}^\mathsf{T}}_{\text{(a) content-to-content}} + \underbrace{\boldsymbol{Q_i^c}\boldsymbol{K_{\delta(i,j)}^r}^\mathsf{T}}_{\text{(b) content-to-position}} + \underbrace{\boldsymbol{K_j^c}\boldsymbol{Q_{\delta(j,i)}^r}^\mathsf{T}}_{\text{(c) position-to-content}} \tag{4}$$

$$\boldsymbol{H_o} = \mathrm{softmax}\left(\frac{\tilde{\boldsymbol{A}}}{\sqrt{3d}}\right)\boldsymbol{V_c}$$

$\tilde{A}_{i,j}$ is the element of attention matrix $\tilde{\boldsymbol{A}}$, representing the attention score from token i to token j. $\boldsymbol{Q_i^c}$ is the i-th row of $\boldsymbol{Q_c}$. $\boldsymbol{K_j^c}$ is the j-th row of $\boldsymbol{K_c}$. $\boldsymbol{K_{\delta(i,j)}^r}$ is the $\delta(i, j)$-th row of $\boldsymbol{K_r}$ with regarding to relative distance $\delta(i, j)$. $\boldsymbol{Q_{\delta(j,i)}^r}$ is the $\delta(j, i)$-th row of $\boldsymbol{Q_r}$ with regarding to relative distance $\delta(j, i)$. Note that we use $\delta(j, i)$ rather than $\delta(i, j)$ here. This is because for a given position i, position-to-content computes the attention weight of the key content at j with respect to the query position at i, thus the relative distance is $\delta(j, i)$. The position-to-content term is calculated as $\boldsymbol{K_j^c}\boldsymbol{Q_{\delta(j,i)}^r}^\mathsf{T}$. The content-to-position term is calculated in a similar way.

Finally, we apply a scaling factor of $\frac{1}{\sqrt{3d}}$ on $\tilde{\boldsymbol{A}}$ which is important for stabilizing model training [1], especially for large-scale PLMs.

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[1] In this sense, our model shares some similarity to Tensor Product Representation [23, 24, 25] where a word is represented using a tensor product of its filler (content) vector and its role (position) vector.

3.1.1 Efficient implementation

For an input sequence of length N, it requires a space complexity of $O(N^2 d)$ [19, 20, 21] to store the relative position embedding for each token. However, taking content-to-position as an example, we note that since $\delta(i, j) \in [0, 2k)$ and thus the embedding of all possible relative positions are always a subset of $K_r \in R^{2k \times d}$, then we can reuse K_r in the attention calculation for all the queries. In experiments, we set the maximum relative distance k to 512 for pre-training. The disentangled attention weights can be computed efficiently using Algorithm 1. Let δ be the relative position matrix according to (3), i.e. $\delta[i, j] = \delta(i, j)$. Instead of allocating a different relative position embedding matrix for each query, we multiply each *query* vector $Q_c[i, :]$ by $K_r^\mathsf{T} \in R^{d \times 2k}$, as in line $3 - 5$. Then, we extract the attention weight using the relative position matrix δ as the index, as in line $6 - 10$. To compute the position-to-content attention score, we calculate $\tilde{A}_{p \to c}[:, j]$, i.e. the column vector of the attention matrix $\tilde{A}_{p \to c}$, by multiplying each *key* vector $K_c[j, :]$ by Q_r^T, as in line $11 - 13$. Finally, we extract the corresponding attention score via the relative position matrix δ as the index, as in line $14 - 18$. In this way, we do not need to allocate memory to store a relative position embedding for each query and thus reduce the space complexity to $O(kd)$ (for storing K_r and Q_r).

Algorithm 1 Disentangled Attention

Input: Hidden state H, relative distance embedding P, relative distance matrix δ. Content projection matrix $W_{k,c}, W_{q,c}, W_{v,c}$, position projection matrix $W_{k,r}, W_{q,r}$.
1: $K_c = HW_{k,c}, Q_c = HW_{q,c}, V_c = HW_{v,c}, K_r = PW_{k,r}, Q_r = PW_{q,r}$
2: $A_{c \to c} = Q_c K_c^\mathsf{T}$
3: **for** $i = 0, ..., N - 1$ **do**
4:     $\tilde{A}_{c \to p}[i, :] = Q_c[i, :]K_r^\mathsf{T}$
5: **end for**
6: **for** $i = 0, ..., N - 1$ **do**
7:     **for** $j = 0, ..., N - 1$ **do**
8:         $A_{c \to p}[i, j] = \tilde{A}_{c \to p}[i, \delta[i, j]]$
9:     **end for**
10: **end for**
11: **for** $j = 0, ..., N - 1$ **do**
12:     $\tilde{A}_{p \to c}[:, j] = K_c[j, :]Q_r^\mathsf{T}$
13: **end for**
14: **for** $j = 0, ..., N - 1$ **do**
15:     **for** $i = 0, ..., N - 1$ **do**
16:         $A_{p \to c}[i, j] = \tilde{A}_{p \to c}[\delta[j, i], j]$
17:     **end for**
18: **end for**
19: $A = \frac{A_{c \to c} + A_{c \to p} + A_{p \to c}}{\sqrt{3d}}$
20: $H_o = \text{softmax}(\frac{A}{\sqrt{3d}})V_c$
Output: H_o

3.2 Enhanced Mask Decoder

During standard BERT pre-training, we feed the final hidden vectors corresponding to the masked tokens to an output softmax over the vocabulary. During fine-tuning, we feed the BERT output into a task-specific decoder, which has one or more task-specific decoding layers, plus a softmax layer if the output is probabilities. To mitigate the mismatch between pre-training and fine-tuning, we treat MLM the same as other downstream tasks, and replace the output softmax of BERT with an Enhanced Mask Decoder (EMD) which contains one or more Transformer layers and a softmax output layer. This makes DeBERTa, which combines BERT and EMD for pre-training, an encoder-decoder model.

The model architecture of the DeBERTa for pre-training is designed to meet several requirements. First, the encoder should be much deeper than the decoder since the former is used for fine-tuning. Second, the number of parameters of the encoder-decoder model needs to be similar to that of BERT so that the pre-training cost is at the same level as BERT. Third, the pre-trained encoder of DeBERTa

should be similar to BERT such that their fine-tuning costs and performances on downstream tasks are comparable.

Take the 12-layer BERT base model as a baseline for comparison. The architecture of DeBERTa for pre-training, used in our experiments, is composed of an encoder that consists of 11 layers of Transformer, a decoder that consists of 2 Transformer layers whose parameters are shared and a softmax output layer. So the model has a similar amount of free parameters as BERT base[2]. After the DeBERTa model is pre-trained, we stack the 11-layer encoder and one decoder layer to recover the standard BERT base architecture for fine-tuning.

In addition, we make a small but important modification on the encoder output vectors before we feed them into the EMD for prediction during MLM pre-training. The authors of BERT [4] propose to not replace all masked tokens with the `[MASK]` token, keeping 10% of them unchanged. Although this is motivated by mitigating the mismatch between fine-tuning and pre-training since `[MASK]` never occurs in the input of downstream tasks, the method suffers from information leaking i.e., predicting a masked token conditioned on the token itself. To address the issue, we replace the encoder output vectors of those masked but unchanged tokens with their corresponding absolute position embedding vectors before we feed them into the decoder for prediction.

4 Experiment

This section evaluates DeBERTa on various NLP tasks.

4.1 Main Results

Following the previous papers on BERT, RoBERTa and XLNet, we report our results using large and base models.

4.1.1 Performance on Large Models

Model	Model Size	CoLA Mcc	QQP F1/Acc	MNLI-m/mm Acc	SST-2 Acc	STS-B Acc	QNLI Acc	RTE Acc	MRPC Acc
BERT$_{large}$	335M	60.6	91.3	86.6/-	93.2	90.0	92.3	70.4	88.0
RoBERTa$_{large}$	355M	68.0	92.2	90.2/90.2	96.4	92.4	93.9	86.6	90.9
XLNet$_{large}$	340M	69.0	92.3	90.8/90.8	**97.0**	92.5	94.9	85.9	90.8
DeBERTa$_{large}$	390M	**69.5**	**92.3**	**91.1/91.1**	96.5	**92.5**	**95.3**	**88.1**	**92.5**

Table 1: Comparison results on the GLUE development set.

We pre-train our large models following the setting of BERT [4], except that we use the BPE vocabulary as [2, 5]. For training data, we use Wikipedia (English Wikipedia dump[3]; 12GB), BookCorpus [26] (6GB), OPENWEBTEXT (public Reddit content [27]; 38GB), and STORIES (a subset of CommonCrawl [28]; 31GB). The total data size after data deduplication[29] is about 78GB. We use 6 DGX-2 machines with 96 V100 GPUs to train the model. A single model trained with 2K batch size and 1M steps takes about 20 days. Refer to Appendix A for the detailed hyperparamters.

We summarize the results on eight GLUE [30] tasks in Table 1, which compares DeBERTa with previous models with around 350M parameters: BERT, RoBERTa, and XLNet. Note that RoBERTa and XLNet use 160G training data while DeBERTa uses 78G training data. RoBERTa and XLNet are trained for 500K steps with 8K samples in a step, which amounts to four billion passes over training samples. We train DeBERTa for one million steps with 2K samples in each step. This amounts to two billion passes of its training samples, approximately half of either RoBERTa or XLNet. Table 1 shows that compared to BERT and RoBERTa, DeBERTa is consistently better across all the tasks. Meanwhile, DeBERTa outperforms XLNet in six out of eight tasks. Particularly, the improvements on MRPC (1.7% over XLNet and 1.6% over RoBERTa), RTE (2.2% over XLNet and 1.5% over

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[2]The disentangled attention layer in DeBERTa has few more free parameters than the self-attention layer in BERT.

[3]https://dumps.wikimedia.org/enwiki/

RoBERTa) and CoLA (0.5% over XLNet and 1.5% over RoBERTa) are significant. Note that MNLI is often used as an indicative task to monitor the progress of pre-training. DeBERTa significantly outperforms all existing models of similar size on MNLI and creates a new state-of-the-art (SOTA).

Model	MNLI-m/mm Acc	SQuAD v1.1 F1/EM	SQuAD v2.0 F1/EM	RACE Acc	ReCoRD F1/EM	SWAG Acc	NER F1
BERT$_{large}$ [4]	86.6/-	90.9/84.1	81.8/79.0	72.0	-	86.6	92.8
RoBERTa$_{large}$ [5]	90.2/90.2	94.6/88.9	89.4/86.5	83.2	90.6/90.0	89.9	93.4
XLNet$_{large}$ [6]	90.8/90.8	95.1/89.7	90.6/87.9	85.4	-	-	-
Megatron$_{336M}$ [29]	89.7/90.0	94.2/88.0	88.1/84.8	83.0	-	-	-
DeBERTa$_{large}$	**91.1/91.1**	**95.5/90.1**	**90.7/88.0**	**86.8**	**91.4/91.0**	**90.8**	**93.8**
Megatron$_{1.3B}$ [29]	90.9/91.0	94.9/89.1	90.2/87.1	87.3	-	-	-
Megatron$_{3.9B}$ [29]	91.4/91.4	95.5/90.0	91.2/88.5	89.5	-	-	-

Table 2: Results on MNLI in/out-domain, SQuAD v1.1, SQuAD v2.0, RACE, ReCoRD, SWAG, CoNLL 2003 NER development set. Note that missing results in literature are signified by "-".

We evaluate DeBERTa on additional benchmarks: (1) Question Answering: SQuAD v1.1 [31], SQuAD v2.0 [32], RACE [33], ReCoRD [34] and SWAG [35]; (2) Natural Language Inference: MNLI [36]; and (3) NER: CoNLL-2003. For comparison, we also include Megatron [29] with three different model sizes: Megatron$_{336M}$, Megatron$_{1.3B}$ and Megatron$_{3.9B}$, which are trained using the same dataset as RoBERTa. Note that Megatron$_{336M}$ has a similar model size as other models mentioned above[4].

We summarize the results in Table 2. Compared to the previous SOTA models with similar sizes, including BERT, RoBERTa, XLNet and Megatron$_{336M}$, DeBERTa consistently outperforms them in all the 7 tasks. Taking RACE as an example, DeBERTa is significantly better than previous SOTA XLNet with an improvement of 1.4% (86.8% vs. 85.4%). Although Megatron$_{1.3B}$ is 3 times larger than DeBERTa, we observe that DeBERTa can still outperform Megatron$_{1.3B}$ in three of the four benchmarks. All the results show the superior performance of DeBERTa in various downstream tasks. We are confident that DeBERTa can perform even better with a larger model size – we leave it to future work.

4.1.2 Performance on Base Models

The setting for base model pre-training is similar to that for large models. The base model structure follows that of the BERT base model, i.e. $L = 12, H = 768, A = 12$. We use 4 DGX-2 with 64 V100 GPUs to train the base model and it takes about 10 days to finish a single pre-training of 1M training steps with batch size 2048. We train DeBERTa with the same 78G text data, and compare it with RoBERTa and XLNet trained using their 160G text data.

We summarize the results in Table 3. Across all three tasks, DeBERTa consistently surpasses RoBERTa and XLNet, with more improvements than that in large models. For example, on the MNLI in-domain setting (MNLI-m), DeBERTa$_{base}$ obtains 1.2% (88.8% vs. 87.6%) over RoBERTa$_{base}$, and 2% (88.8% vs. 86.8%) over XLNet$_{base}$.

Model	MNLI-m/mm (Acc)	SQuAD v1.1 (F1/EM)	SQuAD v2.0 (F1/EM)
RoBERTa$_{base}$ [5]	87.6/-	91.5/84.6	83.7/80.5
XLNet$_{base}$ [6]	86.8/-	-/-	-/80.2
DeBERTa$_{base}$	**88.8/88.5**	**93.1/87.2**	**86.2/83.1**

Table 3: Comparison results on MNLI in/out-domain (m/mm), SQuAD v1.1 and v2.0 development set.

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[4]Although T5 [9] has more parameters (11B), it only reports the test results and it is not comparable with other models.

4.2 Model Analysis

In this section, we first present an ablation study to quantify the relative contributions of different components introduced in DeBERTa. Next, we illustrate the difference in attention patterns between DeBERTa and its counterpart RoBERTa. Last, we study the convergence property to characterize the model training efficiency. We run experiments for analysis using the base model setting where the Wikipedia + Bookcorpus data is used for model pre-training and a model can be pre-trained for 1M steps with batch size 256 in 7 days on a DGX-2 machine with 16 V-100 GPUs.

4.2.1 Ablation study

To verify our experimental setting, we pre-train the RoBERTa base model from scratch. We call the re-pre-trained RoBERTa RoBERTa-ReImp$_{base}$. To investigate the relative contributions of different components in DeBERTa, we design three variations:

- -EMD is the DeBERTa base model without EMD.
- -C2P is the DeBERTa base model without the content-to-position term ((c) in Eq. 4).
- -P2C is the DeBERTa base model without the position-to-content term ((b) in Eq. 4). As XLNet also used relative position bias, this model is close to XLNet plus EMD.

Model	MNLI-m/mm Acc	SQuAD v1.1 F1/EM	SQuAD v2.0 F1/EM	RACE Acc
BERT$_{base}$ [4]	84.3/84.7	88.5/81.0	76.3/73.7	65.0
RoBERTa$_{base}$ [5]	84.7/-	90.6/-	79.7/-	65.6
XLNet$_{base}$ [6]	85.8/85.4	-/-	81.3/78.5	66.7
RoBERTa-ReImp$_{base}$	84.9/85.1	91.1/84.8	79.5/76.0	66.8
DeBERTa$_{base}$	**86.3/86.2**	**92.1/86.1**	**82.5/79.3**	**71.7**
-EMD	86.1/86.1	91.8/85.8	81.3/78.0	70.3
-C2P	85.9/85.7	91.6/85.8	81.3/78.3	69.3
-P2C	86.0/85.8	91.7/85.7	80.8/77.6	69.6
-(EMD+C2P)	85.8/85.9	91.5/85.3	80.3/77.2	68.1
-(EMD+P2C)	85.8/85.8	91.3/85.1	80.2/77.1	68.5

Table 4: Ablation study of the DeBERTa base model.

Table 4 summarizes the results on four benchmark datasets. First, comparing RoBERTa with RoBERTa-ReImp, we observe that they perform similarly across all the four benchmark datasets. Thus, we can confidently treat RoBERTa-ReImp as a solid baseline for comparison. Second, we see that removing any one component in DeBERTa results in a sheer performance drop in all the benchmarks. For instance, removing EMD (-EMD) results in a loss of 1.4% (71.7% vs. 70.3%) on RACE, 0.3% (92.1% vs. 91.8%) on SQuAD v1.1, 1.2% (82.5% vs. 81.3%) on SQuAD v2.0, 0.2% (86.3% vs. 86.1%) and 0.1% (86.2% vs. 86.1%) on MNLI-m/mm, respectively. Similarly, removing either *content-to-position* or *position-to-content* leads to consistent performance drops in all the benchmarks. As expected, removing two components results in even more significant deterioration in performance.
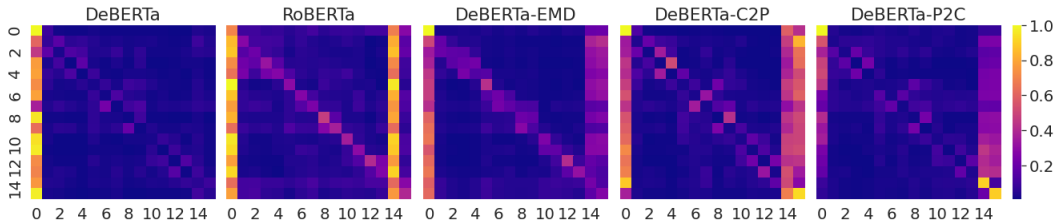

Figure 1: Comparison of attention patterns of the last layer among DeBERTa, RoBERTa and DeBERTa variants (i.e. DeBERTa without EMD, C2P and P2C respectively).

4.2.2 Attention Patterns

To understand why DeBERTa performs differently from RoBERTa, we present their attention patterns in the last self-attention layer in Figure 1, where we also depict the attention patterns of the three DeBERTa variants for comparison. Comparing RoBERTa with DeBERTa, we observe two obvious differences. First, RoBERTa has a clear diagonal line effect for a token to attend to itself, which is not observed in DeBERTa. This could be attributed to the use of EMD, in which the vectors of the masked but unchanged tokens are replaced with their position embeddings. This seems to be verified by examining the attention pattern of DeBERTa-EMD, where the diagonal line effect is brighter than the original DeBERTa. Second, there are vertical strips in the attention patterns of RoBERTa, which are mainly caused by high-frequent functional tokens (e.g., "a", "the", or punctuation). For DeBERTa, the strip appears in the first column, which represents the `[CLS]` token. We conjecture that a dominant emphasis on the `[CLS]` token is desirable for a good pre-trained model since the vector of this token is often used as a contextual representation of the entire input sequence in downstream tasks. We also observe that the vertical strip effect is quite obvious in the patterns of the three DeBERTa variants.

4.2.3 Pre-training Efficiency

To investigate the convergence of model pre-training, we plot the performance of fine-tuned downstream tasks as a function of the number of pre-training steps. As shown in Figure 2, for the RoBERTa-ReImp base model and the DeBERTa base model, we dump a checkpoint every 100K pre-training steps, and then fine-tune the checkpoint on two downstream tasks (MNLI and SQuAD v2.0) and report the accuracy and F1 score, respectively. As a reference, we copy the final model performance of both the original RoBERTa base models [5] and XLNet base models [6] and plot them as flat dot lines. The results show that DeBERTa consistently outperforms RoBERTa-ReImp during the course of pre-training, and converges faster.
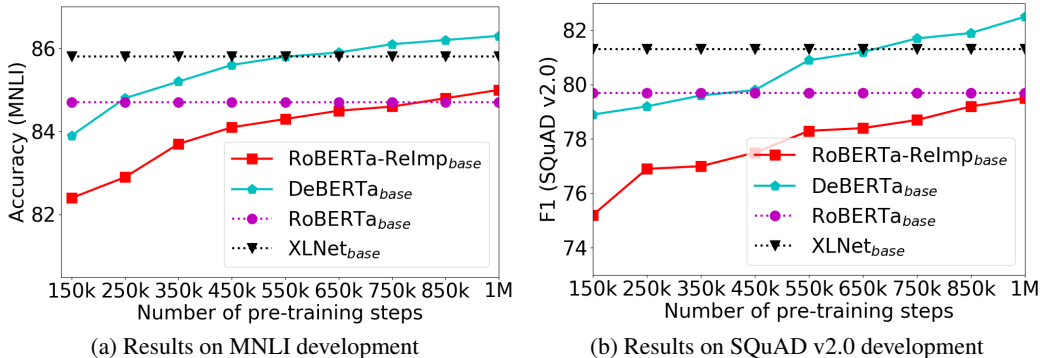

(a) Results on MNLI development	(b) Results on SQuAD v2.0 development

Figure 2: Pre-training performance curve between DeBERTa and its counterparts on the MNLI and SQuAD v2.0 development set.

5 Conclusions

This paper presents two techniques to improve BERT pre-training. The first technique is the disentangled attention mechanism that represents each word using two vectors that encode its content and position, respectively, and computes attention weights among words using disentangled matrices on their contents and relative positions. The second is an enhanced mask decoder that replaces the output softmax layer to predict the masked tokens for MLM pre-training. Using both of these techniques, the new pre-trained language model DeBERTa outperforms RoBERTa and BERT on a wide range of downstream NLP tasks.

This work demonstrates the potential of exploring disentangled word representations for self-attention and the use of task-specific decoders for improving language model pre-training.

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A Appendix

A.1 Dataset

Corpus	Task	#Train	#Dev	#Test	#Label	Metrics
General Language Understanding Evaluation (GLUE)						
CoLA	Acceptability	8.5k	1k	1k	2	Matthews corr
SST	Sentiment	67k	872	1.8k	2	Accuracy
MNLI	NLI	393k	20k	20k	3	Accuracy
RTE	NLI	2.5k	276	3k	2	Accuracy
WNLI	NLI	634	71	146	2	Accuracy
QQP	Paraphrase	364k	40k	391k	2	Accuracy/F1
MRPC	Paraphrase	3.7k	408	1.7k	2	Accuracy/F1
QNLI	QA/NLI	108k	5.7k	5.7k	2	Accuracy
STS-B	Similarity	7k	1.5k	1.4k	1	Pearson/Spearman corr
Question Answering						
SQuAD v1.1	MRC	87.6k	10.5k	9.5k	-	Exact Match (EM)/F1
SQuAD v2.0	MRC	130.3k	11.9k	8.9k	-	Exact Match (EM)/F1
ReCoRD	MRC	101k	10k	10k	-	Exact Match (EM)/F1
RACE	MRC	87,866	4,887	4,934	4	Accuracy
SWAG	Multiple choice	73.5k	20k	20k	4	Accuracy
Token Classification						
CoNLL 2003	NER	14,987	3,466	3,684	8	F1

Table 5: Summary information of the NLP application benchmarks.

- **GLUE**. The General Language Understanding Evaluation (GLUE) benchmark is a collection of nine natural language understanding (NLU) tasks. As shown in Table 5, it includes question answering [31], linguistic acceptability [37], sentiment analysis [38], text similarity [39], paraphrase detection [40], and natural language inference (NLI) [36, 41, 42, 43, 44, 45]. The diversity of the tasks makes GLUE very suitable for evaluating the generalization and robustness of NLU models.

- **ReCoRD** is a commonsense Question Answering dataset. Each example consists of a news article, drawn from CNN and DailyMail, and a Cloze-style question about the article in which one entity is masked out [34].

- **RACE** is a large-scale machine reading comprehension dataset, collected from English examinations in China, which are designed for middle school and high school students [33].

- **SQuAD v1.1/v2.0** is the Stanford Question Answering Dataset (SQuAD) v1.1 and v2.0 [31, 32] are popular machine reading comprehension benchmarks. Their passages come from approximately 500 Wikipedia articles and the questions and answers are obtained by crowdsourcing. The SQuAD v2.0 dataset includes unanswerable questions about the same paragraphs.

- **SWAG** is a large-scale adversarial dataset for the task of grounded commonsense inference, which unifies natural language inference and physically grounded reasoning [35]. SWAG consists of 113k multiple choice questions about grounded situations.

- **CoNLL 2003** is an English dataset consisting of text from a wide variety of sources. It has 4 types of named entity.

A.2 Pre-training Dataset

For DeBERTa pre-training, we use Wikipedia (English Wikipedia dump[6]; 12GB), BookCorpus [26] [7] (6GB), OPENWEBTEXT (public Reddit content [27]; 38GB) and STORIES[8] (a subset of

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[6] https://dumps.wikimedia.org/enwiki/
[7] https://github.com/butsugiri/homemade_bookcorpus
[8] https://github.com/tensorflow/models/tree/master/research/lm_commonsense

CommonCrawl [28]; 31GB). The total data size after data deduplication[29] is about 78GB. For pre-training, we also sample 5% training data as the validation set to monitor the training process.

A.3 Implementation Details

Following RoBERTa [5], we adopted dynamic data batching. We also include span masking[46] as the additional masking strategy with the span size up to three. We list the detailed hyperparameters of pre-training in Table 6. For pre-training, we all use Adam [47] as the optimizer with weight decay [48]. For fine-tuning, even though we can get better and robust results with RAdam[49] on some tasks, e.g. CoLA, RTE and RACE, we all use Adam[47] as the optimizer for a fair comparison. For fine-tuning, we trained each task with a hyper-parameter search procedure, each run will take about 1-2 hours on a DGX-2 node. All the hyperparameters are presented in Table 7. The model selection is based on the performance on the task-specific development sets.

Our code is implemented based on Huggingface Transformers[9], FairSeq[10] and Megatron[29][11].

Hyper-parameter	DeBERTa$_{large}$	DeBERTa$_{base}$	DeBERTa$_{base-ablation}$
Number of Layers	24	12	12
Hidden size	1024	768	768
FNN inner hidden size	4096	3072	3072
Attention Heads	16	12	12
Attention Head size	64	64	64
Dropout	0.1	0.1	0.1
Warmup Steps	10k	10k	10k
Learning Rates	2e-4	2e-4	1e-4
Batch Size	2k	2k	256
Weight Decay	0.01	0.01	0.01
Max Steps	1M	1M	1M
Learning Rate Decay	Linear	Linear	Linear
Adam ϵ	1e-6	1e-6	1e-6
Adam β_1	0.9	0.9	0.9
Adam β_2	0.999	0.999	0.999
Gradient Clipping	1.0	1.0	1.0
Gradient Clipping	1.0	1.0	1.0
Number of DGX-2 nodes	6	4	1
Training Time	20 days	10 days	7 days

Table 6: Hyper-parameters for pre-training DeBERTa.

Hyper-parameter	DeBERTa$_{large}$	DeBERTa$_{base}$
Dropout of task layer	{0,0.1,0.15}	{0,0.1,0.15}
Warmup Steps	{50,100,500,1000}	{50,100,500,1000}
Learning Rates	{5e-6, 8e-6, 9e-6, 1e-5}	{1.5e-5,2e-5, 3e-5, 4e-5}
Batch Size	{16,32,48,64}	{16,32,48,64}
Weight Decay	0.01	0.01
Maximun Training Epochs	10	10
Learning Rate Decay	Linear	Linear
Adam ϵ	1e-6	1e-6
Adam β_1	0.9	0.9
Adam β_2	0.999	0.999
Gradient Clipping	1.0	1.0

Table 7: Hyper-parameters for fine-tuning DeBERTa on down-streaming tasks.

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[9]https://github.com/huggingface/transformers
[10]https://github.com/pytorch/fairseq
[11]https://github.com/NVIDIA/Megatron-LM

A.4 Handling long sequence input

With relative position bias, we choose to truncate the maximum relative distance to k as in (3). Thus in each layer, each token can attend directly to at most $2(k-1)$ tokens and itself. By stacking Transformer layers, each token in the $l-$th layer can attend to at most $(2k-1)l$ tokens implicitly. Taking DeBERTa$_{\text{large}}$ as an example, where $k = 512, L = 24$, in theory, the maximum sequence length that can be handled is 24528. This is a byproduct benefit of our design choice and we found it is beneficial for the RACE task. A comparison of long sequence effect on the RACE task is shown in table 8.

Sequence length	Middle	High	Accuracy
512	88.8	85.0	86.3
768	88.7	86.3	86.8

Table 8: The effect of handling long sequence input for RACE task with DeBERTa

Long sequence handling is an active research area as of late, there are a lot of works built on the Transformer to optimize its performance on long sequence handling[20, 50, 51, 52]. One of the potential future works is to extend DeBERTa to deal with extremely long sequences and compare it with existing approaches.

A.5 Model complexity

With Disentangled Attention, we introduced three additional parameters $\boldsymbol{W_{q,r}}, \boldsymbol{W_{k,r}} \in R^{d \times d}$ and $\boldsymbol{P} \in R^{2k \times d}$. The total increase in parameter is $2L \times d^2 + 2k \times d$. For the large model $(d = 1024, L = 24, k = 512)$, this introduces about $49M$ additional parameters, which is an increment of 13%. For the base model$(d = 768, L = 12, k = 512)$, this introduces about $14M$ additional parameters, which is an increment of 12%.

The additional computational complexity is $O(Nkd)$ due to the calculation of the additional *position-to-content* and *content-to-position* attention scores. Compared with BERT or RoBERTa, this introduces about a 30% increase in computation. Compared with XLNet which also uses relative position embedding, the actual computation cost is about 15%. A further optimization by fusing the attention computation kernel could significantly reduce this additional cost. For EMD, since the decoder in pre-training only reconstructs the masked tokens, it does not introduce additional computation for unmasked tokens. In the situation where 15% tokens are masked and we use only two decoder layers, the additional cost is $0.15 \times 2/L$ which results in an additional computational cost of only 3% for base model$(L = 12)$ and 2% for large model$(L = 24)$ in EMD.

A.6 Attention Patterns

We provide three more examples to illustrate the difference in attention patterns between DeBERTa and RoBERTa.
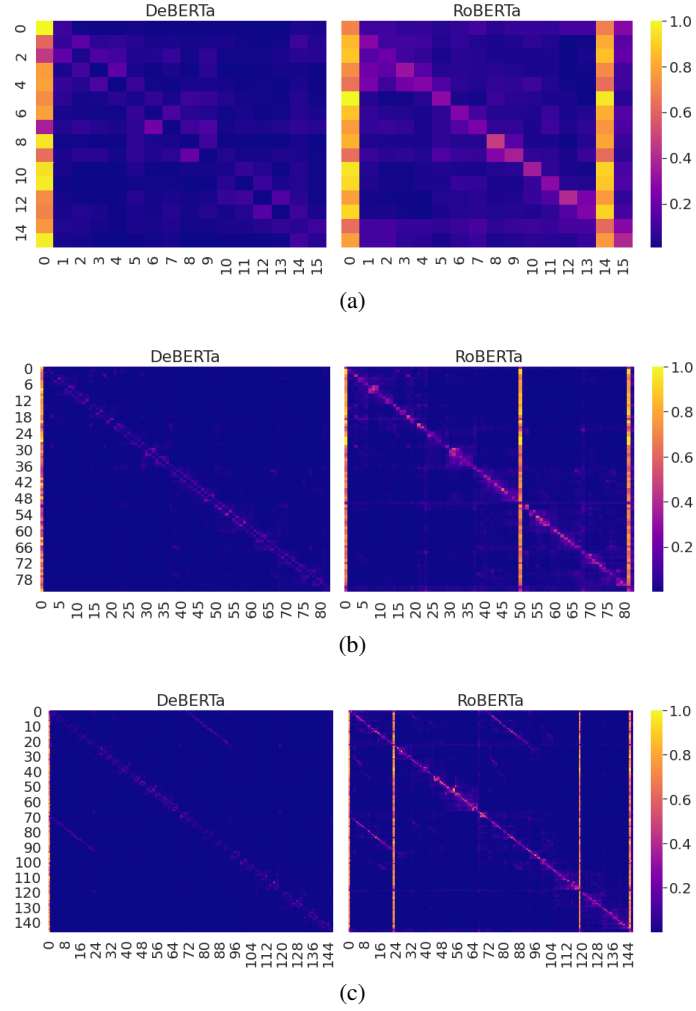
(a)

(b)

(c)

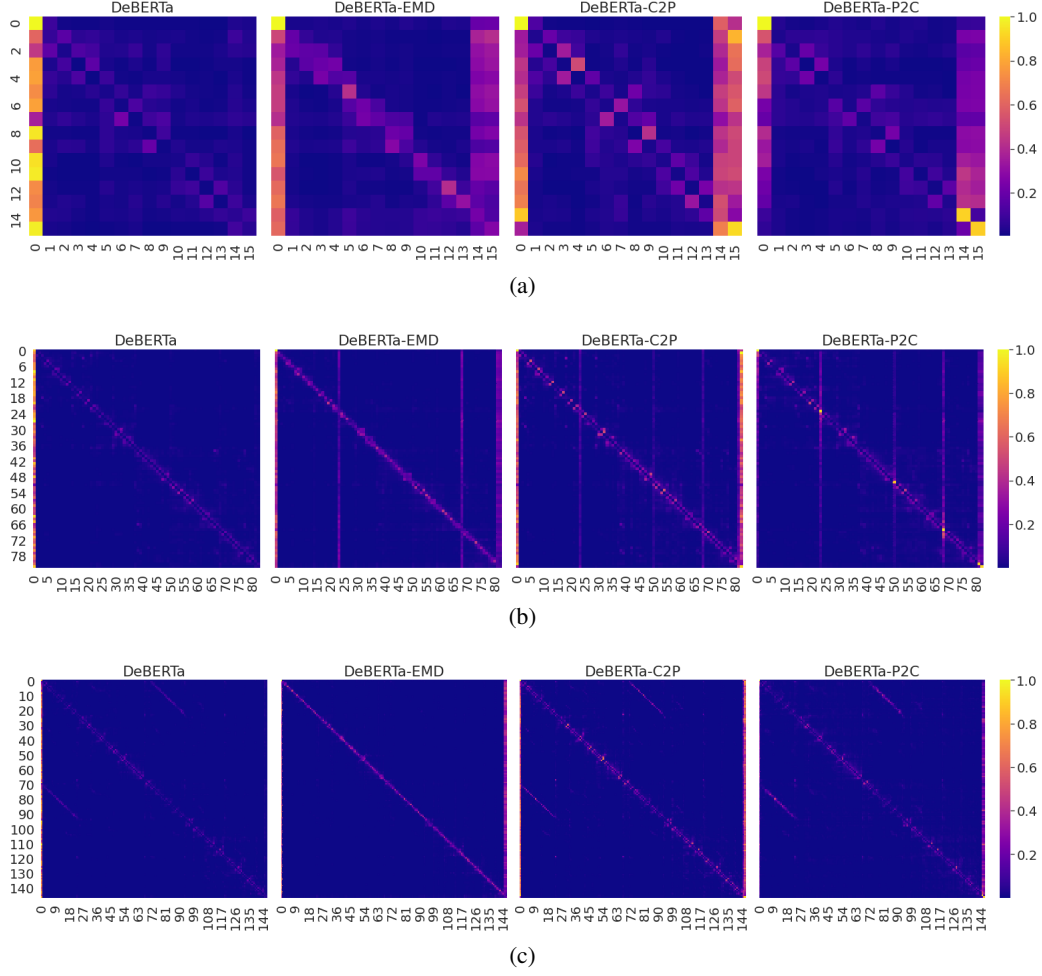Figure 3: Comparison on attention patterns of the last layer between DeBERTa and RoBERTa.

Figure 4: Comparison on attention patterns of last layer between DeBERTa and its variants (i.e. DeBERTa without EMD, C2P and P2C respectively).