Optimized Code Generation for Deep Learning Networks using LATTE and SWIRL

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Abstract

As Deep Neural Networks (DNNs) become more widely used in a variety of applications, the need for performance and portability on a number of different architectures, including CPUs, becomes increasingly important. Traditionally, many DNN frameworks resort to statically-tuned libraries to get performance on specific platforms. This approach is limited by the library performance which can vary greatly across different data sizes and layouts, memory hierarchies and hardware features. Compiler-based methods are getting increased attention because they offer opportunities for performance gains by exploiting data reuse and parallelism, efficient memory access, and vectorization for specific backends with the use of abstraction.

Training DNNs can be challenging, and the Batch Normalization (BN) operator has become a popular technique for accelerating training and making networks more robust. Most DNN frameworks include an optimized implementation of this operator, but the computation efficiency decreases dramatically when this operator does not fit the preoptimized version of library functions.

LATTE is a domain-specific language for DNNs, and SWIRL is a compiler that can
be used with LATTE. This thesis extends the applicability of LATTE/SWIRL by incorporating the BN operator into the LATTE framework and by expanding the optimizations of SWIRL to apply to this operator. Several common compiler techniques, such as scalar replacement, loop interchange, fusion and unrolling, vectorization and parallelization can be applied to this operator for performance enhancements. The optimized BN operator in LATTE/SWIRL is compared to existing frameworks such as TensorFlow, TensorFlow with Intel MKL-DNN, TensorFlow with XLA, PyTorch with MKL-DNN and MXNet with MKL-DNN. The results show that a compiler-based approach for the BN operator can increase performance on CPU architectures.
OPTIMIZED CODE GENERATION FOR DEEP LEARNING NETWORKS USING LATTE AND SWIRL

by

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ABSTRACT

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CHAPTER 1

INTRODUCTION

Deep Neural Networks (DNNs) are currently one of the fastest growing areas in computer science, with wide-ranging applications from speech recognition to genomics. DNNs are a class of models that use a system of interconnected neurons to estimate or approximate functions. Typical DNNs, which consist of multiple stacked layers, require billions of operations for training and inference. These networks are compute intensive, and Graphics Processing Units (GPUs) have been the hardware of choice for many neural networks. However, as the applicability of DNNs has increased, the need for performance on different computing platforms has grown. Due to a variety of factors, including cost and programming complexity, GPUs are not always incorporated into many computing clusters. For example, many scientific applications that employ neural networks are not run on HPC clusters that incorporate GPUs into their architectures. Therefore, there is a demand for performance and portability of DNNs across a variety of architectures and platforms.

Deep Neural Networks are composed of multiple layers, most often including fully-connected, convolutional and activation layers. Many novel layers have been proposed to increase performance and reduce training times. Batch Normalization (BN) is one such layer that standardizes the inputs to other layers. This has the effect of stabilizing the learning process and dramatically reducing the number of epochs required to train deep networks. The use of BN layers in neural networks has quickly become standard due to its effectiveness [2]. Convolutional and fully-connected layers are typically some of the most computationally-intensive sections of the networks. Currently, BN is recognized as one of the most computationally intensive non-convolution layers and is increasingly consuming a larger portion of the execution
time during training for many DNNs [3].

Because of their recent popularity, there are many frameworks that can train and run DNNs. These high-level frameworks, such as TensorFlow [4], Torch [5], Theano [6], Caffe [7], CNTK [8] and MXNet[9], use abstraction to represent neural networks and employ one of three approaches: computation graph engines, layer-specific libraries, and domain-specific languages. Most of these implementations use statically-tuned libraries such as cuDNN for GPUs and Eigen or Intel MKL-DNN for CPUs to achieve performance. cuDNN is a library of highly-tuned implementations for standard routines used in deep neural networks, such as forward and backward convolution, pooling, normalization, and activation layers [10]. Eigen is a high-level C++ library of template headers for linear algebra, matrix and vector operations, geometrical transformations, numerical solvers and related algorithms [11]. Intel MKL-DNN is a performance-enhancing library for accelerating deep learning frameworks on Intel architectures. MKL-DNN implements optimized operators that are common in DNN models, including Convolution, Pooling, Batch Normalization, and Activation. [12].

Due to neural networks being very computationally expensive and involving latency-critical tasks, generating efficient instructions for various platforms is challenging. Preoptimized libraries provide various reliable and fast implementations for linear algebra operations. However, these libraries lack optimization across operators, and the execution of each operation varies dramatically for different data sizes, data layouts, configurations for operators, memory hierarchies and specific hardware features. When a new operator is developed for use in DNNs that does not fit into these preoptimized library functions, the computation efficiency decreases dramatically [13]. Because of these challenges, compiler-based approaches have recently garnered more interest for achieving performance in neural networks. Compiler-based approaches can separate algorithms from schedules, which allows users to experiment with different options for parallelism and data locality on a wide range of platforms. This approach was demonstrated by Halide. Halide is a domain-specific language for image processing that decouples the algorithm definition from the execution strategy and
provides performance portability across different architectures [14].

LATTE is a domain-specific language for DNNs with a graph-like implementation that uses a compiler-based approach for optimization. LATTE provides abstraction to the user through the use of layers. The abstraction hides low-level details such as parallelization, data layout optimizations and code generation from the user. This allows the user to write neural network layers at a high level without architecture-specific optimizations.

SWIRL is a domain-specific compiler for neural networks that can be used with LATTE. SWIRL takes LATTE as input and produces optimized C++ code. Without SWIRL, LATTE uses its own compiler that generates library calls. In contrast, SWIRL uses high-level transformation recipes to generate efficient CPU code. These transformation commands span both data and computation planes. For example, the transformation commands can change the data layout for improved locality, vectorization and parallelization. The layer transformation commands include classical compiler transformations such as tiling, unroll-and-jam and unroll. SWIRL has demonstrated comparable performance with TensorFlow integrated with MKL-DNN on both training and inference for a variety of neural networks, including AlexNet, Overfeat and VGG [15]. Currently, LATTE does not have a Batch Normalization operator. Extending LATTE to include this operator and SWIRL to generate optimizations for this operator will broaden the efficacy and applicability of the LATTE language and the SWIRL compiler for DNNs.

The key contributions of this thesis are:

- An extension of LATTE and SWIRL to include the BN operator and compiler optimizations that can be applied to this operator.

- An application of scalar replacement for reduced memory access and loop interchange and fusion for increased parallelism in the BN LATTE code implementation.
• A transformation recipe for SWIRL to create a SIMD vectorization and parallelization strategy for optimizing BN in LATTE.

• A performance evaluation of the BN operator and of the combined Convolution-BN-ReLU layer on the Intel SkyLake platform, comparing LATTE and SWIRL to TensorFlow, PyTorch and MXNet all integrated with the MKL-DNN library, TensorFlow XLA and native TensorFlow.
CHAPTER 2

BACKGROUND

This section provides a brief description of Batch Normalization and its benefits for training DNNs. The compilation workflow of LATTE and SWIRL is described along with details of how the Batch Normalization Layer is expressed in LATTE.

2.1 Batch Normalization

Advances in deep learning research are largely driven by the introduction of novel hidden layers and architectures. Batch Normalization is a technique introduced in [1] that decreases the training time and increases the robustness of neural networks. Deep neural networks are challenging to train, in part because the input from prior layers can change after weight updates during each training pass. Because the inputs to each layer are affected by the parameters of all preceding layers, this creates an amplifying effect as the network depth increases. Internal covariate shift refers to the variability in the distribution of network activations due to the change in network parameters during training. This variance in the input distribution slows down the training by requiring lower learning rates and careful parameter initialization. Because Batch Normalization reduces the variance in the inputs and the activations in a network, it can allow for higher learning rates during training. Traditionally, higher learning rates are more likely to cause gradients to explode or vanish, which leads to the network getting stuck in local minima during training. Because Batch Normalization reduces the variability in inputs, it prevents small changes to the parameters from amplifying into larger and suboptimal changes in activations for gradients. This reduces the time for convergence.

Covariate shift also makes training difficult with saturating activation functions, such as sigmoid and tanh. Because of the higher variance in inputs, it is more likely that
the input will move into the saturated regime of the nonlinear activation function and slow down convergence. Batch Normalization decreases the variability of the inputs to the nonlinear activation functions, making them more stable and less likely to become saturated, which accelerates the training of the network. The BN transform has been shown to decrease training time and to match performance on many popular DNN models [16]. Because of its proven benefits in training DNNs, its use has become rather ubiquitous in many neural networks today.

While its ability to accelerate the training process in DNNs is not necessarily disputed, the reason for Batch Normalization’s effectiveness has been challenged in recent literature. For instance, the link between the performance gain of BN and the reduction of internal covariate shift has been questioned. In [17], the benefits of BN are demonstrated to be a result of smoothing the objective function. This smoothing creates more predictive gradients and allows for lower learning rates and faster convergence. In [18], the authors argue that the success of Batch Normalization is due to optimizing the length and direction of the parameters separately, which creates a more favorable loss landscape for gradient-based methods.

2.1.1 Batch Normalization Transform

Batch Normalization is achieved through a normalization step that fixes the means and variances of each layer’s inputs. Each dimension of the input data is normalized to a mean of zero and a standard deviation of one. The BN transform is performed on mini-batches since these are used during stochastic gradient training. Therefore, the mean and variation of each input dimension are calculated over a mini batch $B$. Normalizing each input of a layer may change what the layer can represent. To address this, the BN transform includes a pair of parameters, $\gamma$ and $\beta$, which scale and shift the normalized values. These parameters are learned during training. The full Batch Normalization transform algorithm is shown in Figure 2.1.

Equation 1 calculates the mean of a mini-batch $B$ with $m$ inputs. Equation 2
Input: Values of $x$ over a mini-batch: $\mathcal{B} = \{x_1...x_m\}$; Parameters to be learned: $\gamma, \beta$

Output: $\{y_i = BN_{\gamma,\beta}(x_i)\}$

\[
\begin{align*}
\mu_\mathcal{B} &\leftarrow \frac{1}{m} \sum_{i=1}^{m} x_i \quad \text{// mini-batch mean} \\
\sigma^2_\mathcal{B} &\leftarrow \frac{1}{m} \sum_{i=1}^{m} (x_i - \mu_\mathcal{B})^2 \quad \text{// mini-batch variance} \\
\hat{x}_i &\leftarrow \frac{x_i - \mu_\mathcal{B}}{\sqrt{\sigma^2_\mathcal{B} + \epsilon}} \quad \text{// normalize} \\
y_i &\leftarrow \gamma \hat{x}_i + \beta \equiv BN_{\gamma,\beta}(x_i) \quad \text{// scale and shift}
\end{align*}
\]

Figure 2.1: Algorithm for Batch Normalization Transform. Taken from [1]

calculates the variance by subtracting the mean, $\mu_\mathcal{B}$, from each input value and squaring this difference. The variance is also calculated over the mini-batch $\mathcal{B}$. Equation 3 normalizes the inputs to be centered at 0 with a standard deviation of 1. Epsilon is a constant and is added for numerical stability. Equation 4 scales the normalized inputs by $\gamma$ and shifts them by $\beta$.

The BN transform can be added to a network to manipulate any activation. Batch Normalization may be used on the inputs to the layer before or after the activation function in the previous layer. It may be more appropriate after the activation function for S-shaped functions like the hyperbolic tangent and logistic function. For activation functions that can result in non-Gaussian distributions, such as the rectified linear activation function (ReLU), the BN transform is often applied before the activation function [1].

### 2.1.2 Back Propagation

If the BN transform is computed outside the gradient step, the model parameters can blow up and hinder training. Therefore, an important piece of the Batch Normalization technique is allowing the gradient of the loss with respect to the model parameters to account for the normalization. The BN transform is differentiable, and
the gradient of the loss with respect to the different parameters can be computed directly with the chain rule.

\[
\frac{\delta \ell}{\delta \tilde{x}_i} = \frac{\delta \ell}{\delta y_i} \cdot \gamma 
\]

\[
\frac{\delta \ell}{\delta \sigma_B^2} = \sum_{i=1}^{m} \frac{\delta \ell}{\delta \tilde{x}_i} \cdot (x_i - \mu_B) \cdot \frac{-1}{2} \left( \sigma_B^2 + \epsilon \right)^{-3/2} 
\]

\[
\frac{\delta \ell}{\delta \mu_B} = \left( \sum_{i=1}^{m} \frac{\delta \ell}{\delta \tilde{x}_i} \cdot \frac{-1}{\sqrt{\sigma_B^2 + \epsilon}} \right) + \frac{\delta \ell}{\delta \sigma_B^2} \cdot \frac{2(x_i - \mu_B)}{m} 
\]

\[
\frac{\delta \ell}{\delta x_i} = \frac{\delta \ell}{\delta \tilde{x}_i} \cdot \frac{1}{\sqrt{\sigma_B^2 + \epsilon}} + \frac{\delta \ell}{\delta \sigma_B^2} \cdot \frac{2(x_i - \mu_B)}{m} + \frac{\delta \ell}{\delta \mu_B} \cdot \frac{1}{m} 
\]

\[
\frac{\delta \ell}{\delta \gamma} = \sum_{i=1}^{m} \frac{\delta \ell}{\delta y_i} \cdot \tilde{x}_i 
\]

\[
\frac{\delta \ell}{\delta \beta} = \sum_{i=1}^{m} \frac{\delta \ell}{\delta y_i} 
\]

**Figure 2.2: Backpropagation for the Batch Normalization Transform**

The equations shown in Figure 2.2 represent the backward pass of a Batch Normalization training step. The gradient with respect to the normalized inputs, \( \frac{\delta \ell}{\delta \tilde{x}_i} \), is computed in Equation 1. The gradients for \( \gamma \) and \( \beta \) \( \left( \frac{\delta \ell}{\delta \gamma}, \frac{\delta \ell}{\delta \beta} \right) \) are calculated in Equations 5 and 6 respectively.

### 2.2 LATTE and SWIRL

**LATTE** is a domain-specific language for DNNs that provides abstraction for the user to create a neural network, and **SWIRL** uses high-level transformation recipes to generate efficient CPU code. The transformation recipe abstraction allows an expert programmer to explicitly enumerate the transformations that can be applied to each individual layer within **LATTE**. The main **SWIRL** transformations used for the BN operator include tiling, loop unrolling, vectorization and parallelization. Tiling can improve cache locality. Unrolling certain loop iterations by a factor reduces branch penalties and improves register reuse. Vectorization creates intrinsics to be used on SIMD architectures, and parallelization uses OpenMP to parallelize one or more loops.
for increased performance. These optimizations can be tailored for performance on a variety of CPU backends [15].

A DNN is created in LATTE by stacking layers on top of each other, starting with an input layer, adding various hidden layers and ending with a fully-connected layer that applies an activation function. These layers are represented as ensembles of neurons that are connected using mapping functions [19]. LATTE uses an implicit data-flow graph model of the DNN. The nodes represent computations in layers, and the edges are data dependencies between layer inputs and outputs. This data-flow graph is represented by a dictionary of mapping functions, which connects the inputs and outputs of layers. This allows LATTE to store complex graphs without incurring extra memory costs. For training and inference, SWIRL generates kernels for computations as a set of nested for-loops for each forward, backward and weight update pass of each layer.

Figure 2.3 shows a Python code sequence for expressing Batch Normalization in LATTE. The BatchNormLayer function takes three arguments: the network object (net), the input ensemble (input_ensemble) and epsilon. The neurons are created in Line 8 and added to the BN ensemble in Line 10. A mapping function is defined on Lines 13-14 which connects the input_ensemble to the bn_ensemble.

Once the user has defined a neural network in LATTE, the description is then lowered to a standard Python AST. Next, the SWIRL compiler uses transformation recipes on the Python AST. Lastly, the transformed Python AST is translated to C++ code using the ctree package, which is then lowered to optimized x86 machine code using the Intel C++ Compiler. High-quality vector code is also generated via intrinsics rather than relying on compiler directives. See Figure 2.4 for a visual representation of this workflow.
```python
def BatchNormLayer(net, input_ensemble, epsilon=0.001):
    input_channels, input_height, input_width = input_ensemble.shape
    shape = (input_channels, input_height, input_width)
    neurons = np.empty(shape, dtype='object')
    batch_num = net.batch_size

    #Create an ensemble and initialize it
    neurons[:, :, :] = BNNeuron(input_ensemble, batch_num, epsilon)
    bn_ensemble = BNEnsemble(neurons)
    net.add_ensemble(bn_ensemble)

    #Mapping function for add_connections
    def mapping(c, x, y):
        return (range(c), range(x), range(y))
    net.add_connections(input_ensemble, bn_ensemble, mapping)
```

Figure 2.3: Python Code for Batch Normalization Layer in LATTE

Figure 2.4: LATTE/SWIRL Workflow
CHAPTER 3

METHODS

The pseudo code for the Batch Normalization transform, backpropagation and parameter updates is described and shown in Figures 3.1-3.3. These loops can be optimized in a number of ways. Some general optimization techniques, such as scalar replacement combined with loop interchange and loop fusion, were performed on the code directly within LATTE. SWIRL transformation recipes were used for the rest of the optimizations. These are discussed in more detail in this section.

3.1 Batch Normalization in Pseudo Code

The Batch Normalization transform, represented by the equations in Figure 2.1, is expressed by the loop nests in Figure 3.1. Since the transform is applied to mini-batches, the batch dimension $N$ is used. The mean and variance are calculated per feature map or channel dimension $C$, as shown in Lines 6 and 16 of Figure 3.1. The heights and widths of the inputs are $H$ and $W$ respectively. The normalized inputs are referenced as $x_{\text{hat}}$ on Line 26 and are scaled by gamma and shifted by beta on Line 27. The inputs and outputs are $x$ and $y$ respectively.

Figure 3.2 displays the loop nests that express the backpropagation step of Batch Normalization. Because the transform is computed in mini-batches, the gradient is calculated in mini-batches as well. The mini-batch size is $N$. The gradient flowing from the prior layer (\(\nabla y\)) is $\text{grad}$, and $\nabla x$ is $\text{grad}_x$. The gradients for the mini-batch mean and variance (\(\nabla \mu_B, \nabla \sigma_B^2\)) are $\text{dmu}$ and $\text{dvar}$ respectively. The other variables are for ease in computing the gradients. After a series of steps, the gradient with respect to the normalized inputs (referenced as $\text{grad}_x$) is calculated on Line 30 of Figure 3.2.

Figure 3.3 shows the loops nests for the parameter updates of gamma and beta. Sim-
//Calculate the mean per channel dimension
for (n=0; n < N; n++) {
    for(c=0; c < C; c++) {
        for(h=0; h < H; h++) {
            for(w=0; w < W; w++) {
                mean[c] += x[n,c,h,w]
            }
        }
    }
}
for(c=0; c < C; c++) {
    mean[c] = mean[c] / (N * H * W)
}

//Calculate the variance per channel dimension
for (n=0; n < N; n++) {
    for(c=0; c < C; c++) {
        for(h=0; h < H; h++) {
            for(w=0; w < W; w++) {
                var[c] += (x[n,c,h,w]-mean[c])*(x[n,c,h,w]-mean[c])
            }
        }
    }
}
for(c=0; c < C; c++) {
    var[c] = var[c] / (N * H * W)
}

//Apply batch normalization transform
for (n=0; n < N; n++) {
    for(c=0; c < C; c++) {
        for(h=0; h < H; h++) {
            for(w=0; w < W; w++) {
                x_hat[n,c,h,w] = (x[n,c,h,w] - mean[c]) / sqrt(var[c] + epsilon)
                y[n,c,h,w] = gamma[c] * x_hat[n,c,h,w] + beta[c]
            }
        }
    }
}

Figure 3.1: Batch Normalization Forward Pass
for (n=0; n < N; n++) {
    for (c=0; c < C; c++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                divar[c] += grad[n,c,h,w] * gamma[c] * (x[n,c,h,w] - mean[c])
            }
        }
    }
}

dsqrtvar[c] = -1.0 / (var[c] + epsilon) * divar[c]
dvar[c] = 0.5 / sqrt(var[c] + epsilon) * dsqrtvar[c]

for (n=0; n < N; n++) {
    for (c=0; c < C; c++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                dxmu1 = grad[n,c,h,w] * gamma[c] / sqrt(var[c] + epsilon)
dxmu2 = 2.0 * (x[n,c,h,w] - mean[c]) / (N * H * W) * dvar[c]
dx1[n,c,h,w] = dxmu1 + dxmu2
dmu[c] -= dx1[n,c,h,w]
            }
        }
    }
}

for (n=0; n < N; n++) {
    for (c=0; c < C; c++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                dx2[n,c,h,w] = dmu[c] / (N * H * W)
                grad_x[n,c,h,w] = dx1[n,c,h,w] + dx2[n,c,h,w]
            }
        }
    }
}

Figure 3.2: Batch Normalization Back Propagation
ilar to the mean and variance, gamma and beta are calculated per channel dimension $C$. Line 6 computes the gradient for gamma, `grad_gamma`, and Line 7 shows the gradient calculation for beta, `grad_beta`.

```
//Calculate the gradient with respect to the beta and gamma parameters
for (n=0; n < N; n++) {
    for(c=0; c < C; c++) {
        for(h=0; h < H; h++) {
            for(w=0; w < W; w++) {
                grad_gamma[c] += grad[n,c,h,w] * x_hat[n,c,h,w]
                grad_beta[c] += grad[n,c,h,w]
            }
        }
    }
}
```

**Figure 3.3**: Batch Normalization Parameter Update

### 3.2 Scalar Replacement

The goal of scalar replacement is to identify repeated accesses made to the same memory address, either within an iteration or across iterations, and to remove the redundant accesses by keeping the data in registers. Compilers are effective in allocating scalar variables to registers but often fail to do so with array references. Data dependences provide opportunities for reuse of array variables in registers through scalar replacement [13, Chapter 8].

All of the loop nests shown in Figures 3.1-3.3 exhibit data dependences that can be exploited through scalar replacement. For example, in Figure 3.1 the array reference on Line 6 for `mean` has both an output and a true dependence carried by all but the $C$ loop. Line 16 has similar dependences for the array reference to `var` and also an input dependence for the array reference to `mean` carried by all of the loops except $C$. In the last loop nest structure, there are input dependences for `mean` and `var` on line 26 and `gamma` and `beta` on line 27 carried by all of the loops except $C$. Lastly, there is a loop-independent true dependence for `x_hat` on line 27.
Figure 3.4 reflects the changes made for scalar replacement in the Batch Normalization forward pass. To fully exploit the benefits of the scalar replacement, loop interchange of the $C$ and $N$ loops is performed. Because the $C$ loop does not carry any dependences, moving the $N$ loop inside of the $C$ loop allows for more reuse of the values by keeping them in registers during the iterations of the $N$, $H$ and $W$ loops. Since the $C$ loop is now the outermost loop and does not carry any dependences, all of the $C$ loops can be fused. This fusion can increase the level of parallelism and data locality that can be exploited in the compiler transformations applied to this code by SWIRL. It also decreases the loop control overhead. Lastly, a final optimization technique is used for the expensive square root and division operations shown on Line 26 of Figure 3.1. This operation is an input dependence carried in the $C$ loop, which means that this time-consuming calculation needs only to be performed once per $C$ loop iteration rather than for each iteration of every loop. A multiplication operator replaces the more expensive division operator. On Line 21 of Figure 3.4 these calculations are stored in a register and reused on Line 26.

The backward pass exhibits even more opportunities for register reuse. The array reference on Line 6 of Figure 3.2 for $\text{divar}$ has both an output and a true dependence carried by all but the $C$ loop. Input dependences for the array references to $\text{gamma}$ and $\text{mean}$ on Line 6 are carried by all but the $C$ loop as well. Line 8 displays a loop-independent true dependence on the reference to $\text{dsqrtvar}$ and Lines 8-9 have an input dependence on $\text{var}$. In the third loop nest, input dependences are carried by all but the $C$ loop for $\text{gamma}$, $\text{var}$, $\text{mean}$ and $\text{dvar}$. On Line 21, $\text{dmu}$ exhibits the same dependences as $\text{divar}$ does above on Line 6. The last loop nest carries an input dependence on $\text{dmu}$ on Line 29.

Scalar replacements are performed for each of the dependences listed above. As was done in the forward pass, the $C$ and $N$ loops are interchanged and all the $C$ loops are fused. Expensive operations that are input dependences are replaced with scalars in lines 12-14 and 29 of Figure 3.5, which shows the optimized loops nests.
The parameter update loop nest shown in Figure 3.3 is optimized in a similar manner in Figure 3.6. Lines 6 and 7 have output and true and dependences carried by all but the $C$ loop nests for $\text{grad}\_\text{gamma}$ and $\text{grad}\_\text{beta}$ respectively. Additionally, Lines 6 and 7 show loop-independent input dependences for $\text{grad}$. These array references are replaced with scalar variables, and the $C$ and $N$ loops are interchanged. All of the transformations described in Figures 3.4 to 3.6 were expressed within the $\text{LATTE}$ language. The rest of the optimizations were performed using the $\text{SWIRL}$ compiler.
for (c=0; c < C; c++) {
    divar = 0.0
    gamma_temp = gamma[c]
    mean_temp = mean[c]
    var_temp = var[c]

    // Calculate the gradient with respect to the variance on dimension C
    for (n=0; n < N; n++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                divar += grad[n,c,h,w] * gamma_temp * (inputs[n,c,h,w] - mean_temp)
            }
        }
    }
    inver = 1.0 / sqrt(var_temp + epsilon)
    dvar_temp = -0.5 * inver / (var_temp + epsilon) * divar
    inver_dvar = 1.0 / (N * H * W) * dvar_temp

    // Calculate the gradient with respect to the mean on dimension C
    for (n=0; n < N; n++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                dxmu1 = grad[n,c,h,w] * gamma_temp * inver
                dxmu2 = 2.0 * (inputs[n,c,h,w] - mean_temp) * inver_dvar
                dx1_temp = dxmu1 + dxmu2
                dx1[n,c,h,w] = dx1_temp
                dmu += - dx1_temp
            }
        }
    }

    // Calculate the gradient with respect to the inputs
    inver_dmu = 1 / (N * C * W) * dmu
    for (n=0; n < N; n++) {
        for (h=0; h < H; h++) {
            for (w=0; w < W; w++) {
                dx2[n,c,h,w] = dmu / N
                grad_x[n,c,h,w] = dx1[n,c,h,w] + inver_dmu
            }
        }
    }
}

Figure 3.5: Batch Normalization Back Propagation with Scalar Replacement Optimizations
//Calculate the gradient with respect to the beta and gamma parameters
for(c=0; c < C; c++) {
    grad_gamma_temp = 0.0
    grad_beta_temp = 0.0
    for (n=0; n < N; n++) {
        for(h=0; h < H; h++) {
            for(w=0; w < W; w++) {
                grad_temp = grad[n,c,h,w]
                grad_gamma_temp += grad_temp * x_hat[n,c,h,w]
                grad_beta_temp += grad_temp
            }
        }
    }
    grad_gamma[c] = grad_gamma_temp
    grad_beta[c] = grad_beta_temp
}

Figure 3.6: Batch Normalization Parameter Update with Scalar Replacement Optimizations

3.3 Transformation Recipes

SWIRL uses the concept of transformation recipes to generate high-quality code and to customize optimizations to the target hardware. A recipe consists of a set of commands that include both data and code transformations. For a more detailed explanation of these commands, the reader is directed to [15].

The transformation recipe used for the BatchNormLayer forward pass implementation is shown in Figure 3.7. The feature map dimensions of the input and output are tiled by the SIMD vector length for the given platform to aid vectorization (Lines 4-5). The loops corresponding to the batch dimension are specified for concurrent execution via the parallelize command in Lines 6 and 8. During code generation these loops are annotated with OpenMP pragmas. The inner-most loop of the 4-D loop nest is vectorized and unrolled by a factor that allows for the data locality to be fully exploited by the SIMD instructions and to reduce loop overhead (Lines 10-11).

A sample transformation recipe for the backward and weight update passes used in the BatchNormLayer is illustrated in Figure 3.8. The transformations are similar to the ones described above for the forward pass. During training, grad_inputs represents
if "value" in input_ensemble.tiling_info:
    tiled_dims = input_ensemble.tiling_info["value"]
    for dim, factor in tiled_dims:
        bn_ens.tile('inputs', dim=dim, factor=factor)
        bn_ens.tile('value', dim=dim, factor=factor)
        bn_ens.parallelize(phase="forward", loop_var="_neuron_index_1_outer")
else:
    bn_ens.parallelize(phase="forward", loop_var="_neuron_index_1")

bn_ens.vectorize(phase="forward", loop_var="_neuron_index_3", factor=latte.config.SIMDWIDTH)
bn_ens.unroll(phase="forward", loop_var="_neuron_index_3", factor=4, unroll_type=0)

**Figure 3.7:** An example transformation recipe for the forward pass of the Batch Normalization layer created in Figure 2.3.

the input gradients and `grad` represents the output gradients of the current ensemble layer. The feature map dimensions of the `grad_input` and `grad` are tiled by the SIMD vector length for the given platform to aid vectorization (Lines 4-5). The loops corresponding to the batch dimension are specified for parallelization for both the backward and weight update loops via the parallelize command in Lines 6 and 8. The inner-most loop is vectorized and unrolled as shown in Lines 10-11.

Applying the transformation recipes in Figures 3.7 and 3.8 to the Batch Normalization layer in *LATTE* generates optimized code for the Intel SkyLake platform. The final C++ code generated by the *SWIRL* compiler for the forward pass is shown in Figure 3.9.
if "grad" in input_ensemble.tiling_info:
    tiled_dims = input_ensemble.tiling_info["grad"]
    for dim, factor in tiled_dims:
        bn_ens.tile('grad_inputs', dim=dim, factor=factor)
        bn_ens.tile('grad', dim=dim, factor=factor)
        bn_ens.parallelize(phase="backward", loop_var="_neuron_index_1_outer")
else:
    bn_ens.parallelize(phase="backward", loop_var="_neuron_index_1")

bn_ens.vectorize(phase="backward", loop_var="_neuron_index_3", factor=latte.config.SIMDWIDTH)
bn_ens.unroll(phase="backward", loop_var="_neuron_index_3", factor=4, unroll_type=0)

Figure 3.8: An example transformation recipe for the backward and weight update pass of the Batch Normalization layer created in Figure 2.3
for (int neuron_index_0 = 0; neuron_index_0 < 2; neuron_index_0 += 1) {
    double var_t = 0.0;
    for (int neuron_index_2 = 0; neuron_index_2 < 64; neuron_index_2 += 1) {
        var_t += _mm512_reduce_add_ps(var_temp_3, mean_t2);
        mean_t2 += _mm512_reduce_add_ps(mean_temp_3, _mm512_set1_ps(8192));
        mean_temp_3 = _mm512_add_ps(mean_temp_3, _mm512_load_ps(&ensemble2inputs[_neuron_index_0][_neuron_index_2][(_neuron_index_3 + 48)]));
        var_temp_3 = _mm512_fmadd_ps(di_ff_3, di_ff_2, var_temp_2);
        var_temp_2 = _mm512_fmadd_ps(di_ff_2, di_ff_1, var_temp_1);
        var_temp_1 = _mm512_fmadd_ps(di_ff_1, di_ff_0, var_temp_0);
        var_temp_0 = _mm512_fmadd_ps(di_ff_0, di_ff_3, var_temp_3);
    }
}

for (int neuron_index_1 = 0; neuron_index_1 < 2; neuron_index_1 += 1) {
    double mean_t1 = 0.0;
    for (int neuron_index_0 = 0; neuron_index_0 < 2; neuron_index_0 += 1) {
        mean_t1 += mm512_reduce_add_ps(mean_temp_0, mm512_set1_ps(0.0));
        mean_temp_0 = _mm512_add_ps(mean_temp_0, _mm512_load_ps(&inputs[_neuron_index_0][_neuron_index_1][_neuron_index_2][_neuron_index_3]));
    }
    mean_t1 = _mm512_reduce_add_ps(mean_temp_0, _mm512_set1_ps(8192));
    mean_temp_0 = _mm512_add_ps(mean_temp_0, _mm512_load_ps(&inputs[_neuron_index_0][_neuron_index_1][_neuron_index_2][_neuron_index_3]));
    mean_t1 = _mm512_reduce_add_ps(mean_temp_0, _mm512_set1_ps(8192));
    mean_temp_0 = _mm512_add_ps(mean_temp_0, _mm512_load_ps(&inputs[_neuron_index_0][_neuron_index_1][_neuron_index_2][_neuron_index_3]));
}

mean_t = _mm512_div_ps(mean_t1, _mm512_set1_ps(2.0));

for (int neuron_index_1 = 0; neuron_index_1 < 2; neuron_index_1 += 1) {
    for (int neuron_index_2 = 0; neuron_index_2 < 64; neuron_index_2 += 1) {
        mean_t = _mm512_fmadd_ps(gamma_temp, x_hat_2, beta_temp);
        x_hat_2 = _mm512_fmadd_ps(gamma_temp, x_hat_1, beta_temp);
        x_hat_1 = _mm512_fmadd_ps(gamma_temp, x_hat_0, beta_temp);
        x_hat_0 = _mm512_fmadd_ps(gamma_temp, _mm512_sub_ps(_mm512_load_ps(&inputs[_neuron_index_0][_neuron_index_1][_neuron_index_2][_neuron_index_3]), mean_t), 1.0 / sqrt(var_t + 0.001));
    }
}

Figure 3.7: Generated C++ code from the SWIRL transformation recipe shown in Figure 3.7
CHAPTER 4
RESULTS

The performance results of using LATTE and SWIRL for the Batch Normalization layer compared to other state-of-the-art frameworks is described in this section. The results were generated on an Intel Skylake platform with AVX-512 support. The frameworks used for comparison include TensorFlow release version 1.11.0, TensorFlow release version 2.0.0 configured with Intel Math Kernel Library (MKL), TensorFlow release version 2.0.0 using XLA, MXNet version 1.5.1 with MKL, and Pytorch version 1.4.0+cpu with MKL.

4.1 Hardware Platform and Environment
The hardware platform used is a high performance server class dual socket Intel Xeon Gold 6130 SkyLake processor with $2 \times 32$ 2.1 Ghz (max 3.7 Ghz) turbo-enabled cores. This is an AVX-512 platform with 512-bit vector support. The processor has 98 GB of DDR4-2666 memory, with 32KB of L1 cache, 1MB of L2 cache and 22MB of L3 Cache. The code is generated via the Intel C++ Compiler (ICC) v18.0.1.163 with "-O3 -qopenmp -xCORE-AVX512" flags and NUM_OMP_THREADS=32.

4.2 Performance Comparison of Batch Normalization
The performance of the LATTE/SWIRL implementation of the BN operator used for training is compared with the five frameworks described above: TensorFlow, TensorFlow with MKL-DNN, TensorFlow using XLA, MXNet with MKL-DNN, and PyTorch with MKL-DNN. The training step involves a forward, backward and weight update pass. The testing is carried out on ten different input sizes. The dimensions of these input layers are representative of the sizes found in GoogleNet[20], VGGNet[21] and ResNet[22] architectures. The results are displayed in Figure 4.1, comparing
the number of images per second each implementation can process with a batch size of 64 and the image size listed as C, H/W where C represents the number of channels and H/W represent the height and width dimensions respectively. LATTE/SWIRL outperforms all of the other frameworks. On average, LATTE/SWIRL has 2x greater throughput than MXNet with MKL-DNN, 4x more throughput than PyTorch with MKL-DNN, 5x greater throughput than TensorFlow using XLA, 6x more throughput than TensorFlow with MKL-DNN and 100x greater throughput than TensorFlow.

Figure 4.1: Performance results and breakdown for the Batch Normalization training step. Demonstrates the effects of different optimizations towards overall performance for the LATTE/SWIRL implementation.
The code optimizations shown in Figures 3.4-3.5 and the transformation recipes shown in Figures 3.6-3.7 incorporate several compiler optimizations, including scalar replacement, loop unrolling, SIMD vectorization and parallelization. The individual effects of each optimization on the LATTE/SWIRL implementation are included in the performance results of the Batch Normalization operator in Figure 4.1. The breakdown is shown only for LATTE/SWIRL results, with the blue area representing the baseline performance, the orange area displaying the performance with scalar replacement included, the grey area showing the performance with the loop unrolling and SIMD vectorization in addition to scalar replacement, and the gold area exhibiting the performance with all of the optimizations, including parallelization. Scalar replacement achieves an increase between 4% to 18%, with an average of 15%. The loop unrolling and SIMD vectorization boost performance between 2% and 11%, with an average of 9% and parallelization provides the greatest benefit by improving performance from 69% to 93%, with an average increase of 77%. The baseline performance of LATTE/SWIRL is roughly 2x better than the TensorFlow implementation without MKL-DNN.

4.3 Performance Comparison for Conv-BN-ReLU Layer

A comparison is presented for a training layer composed of Convolution, Batch Normalization and ReLU activation operators. This layer configuration is commonly found in many deep learning architectures and is suggested in [1]. The layer dimensions are representative of layers found in the GoogleNet[20], VGGNet [21] and ResNet[22] architectures. The results can be seen in Figures 4.2 and 4.3. Figure 4.2 shows the images per second that each implementation can process with batch sizes of either 64 or 128 with varying image sizes. C represents the number of in-channels for the convolution layer, K represents the out-channels for the Convolution operator and hence the number of channels for the BN and ReLU operators, and H/W are the height and width respectively of the image for all of the operators. Figure 4.3 displays a time comparison for computing the Conv-BN-ReLU layer. This comparison is presented so a breakout of the performance of each operator can be observed. Note that the results from the native TensorFlow implementation are not included so the
graph would not be distorted, and the information from the other frameworks could be seen more easily.

LATTE/SWIRL outperforms the other implementations on most of the tests. For those dimensions where H/W is small (\(<=14\)), the Convolution operator in LATTE/SWIRL is not as efficient as the Convolution operator in TensorFlow with MKL and MXNet with MKL. The largest performance gains by LATTE/SWIRL are observed for the test dimensions where H/W is large (i.e. 112). This demonstrates that the loop unrolling and SIMD vectorization in LATTE/SWIRL on the inner dimensions can be more fully exploited when the inner dimensions are large. For all of the test sizes, LATTE/SWIRL was more efficient than the other frameworks for the BN and ReLU activation operators.
Figure 4.2: Performance results in images per second for the Convolution-Batch Normalization-ReLU layer training step.
Figure 4.3: Performance results in time (ms) for the Convolution-Batch Normalization-ReLU layer training step. Displays the time spent on each operator.
CHAPTER 5

RELATED WORK

Compiler-based approaches are gaining more interest in the quest for performance and portability of DNNs, which means that more compiler-based frameworks are being introduced. TensorFlow’s XLA [23] compiles neural networks for CPUs, GPUs and accelerators. It lowers nodes into primitive linear algebra operations and then calls a backend-specific library for different architectures (such as Eigen for CPUs, or cuDNN for GPUs) to perform the bulk of the computation. It aims to provide backend flexibility for TensorFlow.

TVM [24] is a compiler that exposes graph-level and operator-level optimizations. It lowers the traditional neural network dataflow graph into a two-phase internal representation (IR). The high-level IR is optimized independent of the target architecture. Then it lowers nodes into a low-level Halide-based internal representation wherein loop-based optimizations can be performed. Halide is used to generate LLVM or CUDA/Metal/OpenCL source code. TVM relies on an inventory of optimization recipes. When an input layer’s parameters match one of those in the inventory, the optimizations are invoked. Currently, TVM does not have support for training operations, only inference.

Glow [25] is a machine learning compiler for heterogeneous hardware. Similar to TVM, it uses a two-phase IR. The high-level internal representation allows the optimizer to perform domain-specific optimizations. The lower-level IR permits the compiler to perform memory-related optimizations. Glow then calls optimized linear algebra libraries in the lower-level IR. It is similar to TVM in that uses a computational graph engine for the high-level representation, and it also aims to support
multiple backends. Glow also focuses on inference operators with the intent to focus on training operations in the future.
CHAPTER 6

CONCLUSIONS

Compiler-based approaches have proven to be an effective way to increase portability of DNNs through abstraction while also achieving performance on a variety of architectures. This research project involved extending the LATTE language and the SWIRL compiler to implement the Batch Normalization operator. This work increases the applicability of LATTE/SWIRL for modern DNNs and demonstrates the effectiveness of using compiler-based approaches as compared to other methods, such as statically-tuned libraries. Performance evaluations of this extension were tested at both the operator level and layer level. These tests were run on an Intel SkyLake platform using a variety of input sizes that are found in common network architectures. Performance gains were observed for all of the comparisons at the operator level and most of the tests at the layer level. These results demonstrate the effectiveness of a compiler-based approach for achieving performance and portability for modern neural networks.
REFERENCES


