

TModel: Modeling Texture Memory and Mobile GPU Performance to Accelerate DNN Computations

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Abstract

The demand for Deep Neural Network (DNN) execution (including both inference and training) on mobile system-on-a-chip (SoCs) has surged, driven by factors like the need for real-time latency, privacy, and reducing vendors' costs. Mainstream mobile GPUs (e.g., Qualcomm Adreno GPUs) usually have a 2.5D L1 texture cache that offers throughput superior to that of on-chip memory. However, to date, there is limited understanding of the performance features of such a 2.5D cache, which limits the optimization potential. This paper introduces TModel, a framework with three components: 1) a set of micro-benchmarks and a novel performance assessment methodology to characterize a non-well-documented architecture with 2D memory, 2) a complete analytical performance model configurable for different data access pattern(s), tiling size(s), and other GPU execution parameters for a given operator (and associated size and shape), and 3) a compilation framework incorporating this model and generating optimized code with low overhead. TModel is validated both on a set of DNN kernels and for training complete models on mobile GPU, and compared against both

popular mobile DNN frameworks and another GPU performance model. Evaluation results demonstrate that TModel outperforms all baselines, achieving $1.48 - 3.61\times$ speedup on individual kernels and $1.83 - 66.1\times$ speedup for end-to-end on-device training with only $0.25\% - 18.5\%$ the tuning cost of the baselines.

CCS Concepts

• **Computing methodologies** → **Neural networks**; **Parallel computing methodologies**; • **General and reference** → **Performance**; • **Human-centered computing** → *Ubiquitous and mobile computing systems and tools*; • **Software and its engineering** → *Source code generation*.

Keywords

Texture Memory, Performance Modeling, Architecture Profiling, Automatic Code Generation, On-device Training

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1 Introduction

Mobile system-on-a-chip (SoC) has become an indispensable platform for executing powerful compute- and memory-intensive applications. Prominent among these are deep learning (DL) neural network execution tasks [10, 30] that

support applications related to speech, image, and text recognition, or even Generative AI. Execution on a mobile device itself, as compared to the mobile device serving as a front-end to cloud-based servers, has several advantages like service continuity or lower latency of applications with no or limited internet availability, and better privacy, among others. However, execution on mobile devices is challenging given the constantly increasing size of DL models and the limited resources of mobile devices.

Performance models, especially those related to memory/cache, have long played a crucial role in guiding performance tuning of applications [8, 24, 69, 77, 78, 81]. Modeling on-chip cache behaviors has been extensively explored in both sequential [14, 25] and parallel processors [3, 4, 12, 32, 33, 50, 53, 62, 63, 66]. A prevalent approach involves extracting memory traces and calculating the *data reuse distance* metric. In developing performance models for mobile processors, there are significant new considerations. For example, the latest Adreno and Mali GPUs are equipped with specialized L1 texture cache, offering notable performance advantages over other on-chip memory types (e.g., local and constant memory [39]). The existing performance modeling approaches, including the notion of data reuse distance, do not directly apply to the texture cache, which is designed for two-dimensional spatial locality. An additional complication is that the texture representation (i.e., the storage mapping in memory) is patented and obscure to developers.

Based on this motivation, this paper proposes TMMoel, the first performance model and associated compiler designed to address the following question: *“how to analytically choose between different implementations of a single NN operator for a given input and output shape – specifically, making selections considering data access pattern(s), tile sizes, and other GPU execution parameters – to obtain the best performance”*. This performance model needs to capture (i) the spatial locality of a single thread with texture memory, (ii) the cross-thread spatial and temporal locality, and (iii) the interplay with other GPU factors like occupancy.

The design and implementation of TMMoel comprises several steps. First, we develop a set of micro-benchmarks to demystify the black-box micro-architecture and introduce novel approaches to characterize performance. One of the key ideas we introduce is *cross-block stride*, designed to capture spatial locality in 2D memory. The insights from micro-benchmarks lead to a performance model, integrating the factors listed above into one unified cost expression. Furthermore, the performance model is incorporated into a compiler that can generate efficient code efficiently (i.e., without high performance tuning costs that many other systems require).

Overall, this paper makes the following contributions:

- We introduce a series of micro-benchmarks to understand several aspects of the target architecture – including the

impact of spatial and temporal locality of the texture memory; the work-group size, warp shape, and the occupancy, on the performance. The methodology and the set of micro-benchmarks is designed to be applicable to other architectures as well.

- We propose an automatic way to predict the latency for a given data access pattern on 2D texture memory space.
- We introduce a full performance model, capturing the impacts of memory and occupancy across threads in a warp, warps in a work group, and the full kernel.
- We develop a complete system prototype for code optimization for both kernels and full DL model training on a mobile architecture.

TMMoel is extensively evaluated on both individual kernels and end-to-end on-device DNN training by comparing with four state-of-the-art mainstream product-level mobile DNN frameworks and their associated performance modeling/tuning methods (MNN [30], CLML [1], TFLite [2], and TVM [9]) and a more general-purpose GPU performance modeling method (PPT-GPU [3]). The evaluation results demonstrate that TMMoel achieves an average speedup of $1.48 - 3.61\times$ on individual kernels and a speedup of $1.83 - 66.1\times$ for end-to-end DNN on-device training. TMMoel achieves similar performance to the brute-force search (with an average variation of 1%), with only 0.25% - 18.5% tuning cost vs the baseline frameworks. TMMoel is also characterized by high prediction accuracy and good portability.

2 Background

This section provides an overview of mobile GPU architectures, starting with details of the OpenCL programming model, which is a popular way to program these GPUs. To provide additional context for the entire paper, this section also compares OpenCL with CUDA terminology.

OpenCL Programming Model. OpenCL, like CUDA, employs the SIMT (Single Instruction Multiple Threads) execution model, where a kernel specifies the computation logic for one thread. The data index space associated with this execution is divided into *work groups* (i.e., *thread blocks* in CUDA) based on the specified *work-group size*. Each work group is assigned to a *shader core* (equivalent to a *streaming multiprocessor* in CUDA). A work group consists of multiple *work items*, which are software *threads* mapped to hardware threads for execution on ALUs within the shader core. The fundamental execution unit in a shader core is the *warp*, a group of hardware threads that execute the same instruction simultaneously. Consequently, the work items within a work group are partitioned into warps for execution. The launch order of warps or work groups is not defined, so programs should not assume any specific execution sequence across them [55].

Mobile GPU Architecture and 2D Texture Cache. Fig. 1 shows a high-level overview¹ of the mainstream Adreno mobile GPU architecture by Qualcomm [55]. Different generations of Adreno GPUs or other mobile GPUs (e.g. Mali series) may have different implementations with respect to parameters such as the number of shader cores (#SP) or the capacity of different caches, and other factors [5, 39, 55].

2D texture memory. System memory supports two types of data storage: *buffer* and *image*. The former stores data in 1-dimension continuously as in CPU memory and reads and writes data via L2 cache. The latter allows data storage as images in 1-/2-/3-dimensional entities called *image objects*, with each data element (called *pixel*) capable of storing four channels (RGBA). Usually, mobile GPUs accelerate data reads of *image objects* by a special two-dimensional or 2D L1 cache hardware called *texture cache*². Although the read-only L1 texture cache is small (usually 1KB) and shared among all threads on a shader core, it offers much higher reading throughput than other memory (e.g., 2× over L2 and 7× over main memory [39, 55]). Exploiting 2D data locality turns out to be critical for performance optimization on mobile GPUs. **Differences between Mobile and Desktop GPUs.** Mobile and desktop GPUs differ significantly in their memory hierarchy and memory access speed. On desktop GPUs, the access latency of shared memory (or L1 cache) is 1.47× to 6.87× faster than that of the texture cache – as observed across multiple generations of GPU microarchitectures [18, 35, 41, 44, 51, 70]. In contrast, on mobile GPUs, the texture cache offers twice the throughput of shared memory [39].

On Nvidia GPUs, achieving high memory bandwidth primarily depends on coalescing accesses and minimizing bank conflicts from shared memory. In contrast, mobile GPUs benefit more from vectorized memory access and coalesced accesses are not effective [55]. To further explain the difference, coalesced accesses involve threads in a warp collectively accessing contiguous, aligned addresses, whereas vectorized access allows a single thread to load/store multiple consecutive elements in one operation [55, 71]. Moreover, efficiently leveraging the on-chip texture L1 cache, with its limited 1KB capacity, is crucial for high performance.³

3 Demystify a System with 2D Cache

Understanding texture cache in-depth is necessary in order to build an analytical performance model for applications on a system using such a cache. However, this is challenging because of two key design features (*texture representation* and *cache organization*) of mobile texture cache [15, 22, 26]

¹Graphics-only hardware modules are omitted because this paper focuses on general-purpose computing on mobile GPUs.

²Also referred to as 2.5D as each point in the 2D space is a vector.

³In comparison, shared memory/L1 size on latest server GPUs is 256 KB [52].

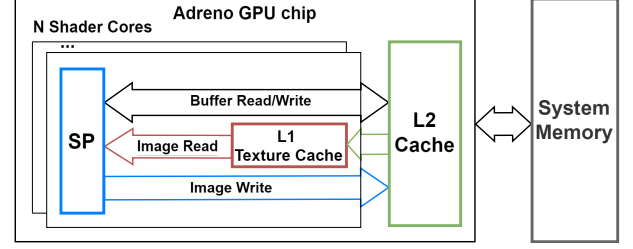


Figure 1: High-level overview of Adreno GPUs.

are patented and obscure to end-users [55]. Texture representation refers to the two-dimensional texture objects' data storage in memory. This further includes the *data unit* (e.g., 2D data block), their *organization* (e.g., hierarchical blocked storage), and the *data storage order* in memory within and among the data units (e.g., row, column, zigzag, or Hilbert order). Cache organization defines the data read between the 2D texture cache and the main memory including data mapping, cache line replacement/eviction, and other features. Previous methods [44, 70] focused solely on using 1D image data (Image1D) to analyze the texture cache organization, leaving the analysis of texture representation, 2D data locality, and their impact on performance unclear. With Image2D outperforming Image1D by a factor of 2 in terms of throughput, it is clearly important to focus on the former. Our work introduces a general empirical approach for studying the relationship between data access patterns and the resulting memory performance. This is accomplished in two steps: considering texture representation and 2D data locality for a single thread (Section 3.1), and extending this study by considering GPU's SIMT parallel programming (Section 3.2).

3.1 Single-Thread Performance with 2D Cache

This empirical study leverages a set of micro-benchmarks involving random accesses to unveil the relationship between data access patterns and 2D texture object memory latency for a single thread. As noted above, the goal is to achieve this without necessarily knowing the concrete texture representation and 2D cache organization. The outcome of this study is a *machine learning regression model* that can predict the expected memory latency for any given data access pattern. Figure 2 illustrates the overview of this study which consists of four main steps outlined below. Constrained by space, please find a complete algorithm from our anonymous Supplementary Material (Sec. I) online⁴.

Step I: Offline multinomial distribution-based random strides generation. Inspired by pointer-chasing benchmarking [65],

⁴<https://drive.google.com/file/d/1UYV8sxYRva1EhSad3qp6U8Od-YWb9kq>

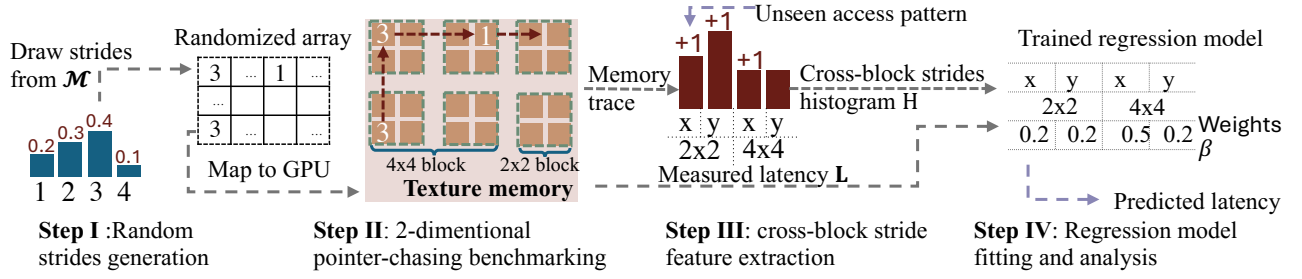


Figure 2: Overview of the empirical study for single-thread data access pattern and 2D texture cache performance. **Step I** generates a random stride list of (1, 2, 3, 4) with a possibility of (0.2, 0.3, 0.4, 0.1) in a multinomial distribution. **Step II** conducts a 2D pointer-chasing benchmarking using random strides and generates a memory trace. **Step III** extracts cross-block strides and creates a histogram as a regression model feature. This example assumes 2 block shapes, 2×2 and 4×4 , counts cross-block strides along both x- and y-directions for both cases, and creates a cross-block strides histogram w/ four bars. **Step IV** runs multiple repetitions of the first three steps and trains a regression model w/ multiple cross-block stride histograms and the corresponding measured latency, which can predict the access latency for arbitrary data access patterns later.

this step generates random access indices offline, thus eliminating the overhead of online index generation and memory trace collection, and mitigating their impact on benchmarking accuracy. Unlike existing efforts [39, 44, 65, 70] that generate 1D uniform strides, TMMModel is novel in generating 2D random strides (along horizontal and vertical directions) to study the 2D texture object access. This step takes a list of possible strides (S) as input and constructs a multinomial distribution $\mathcal{M}$ for S . Each benchmark execution assumes a different multinomial distribution $\mathcal{M}$ to ensure randomness.

Step II: On-device 2D Random-strided pointer-chasing benchmarking. After generating random strides, TMMModel employs a newly designed micro-benchmark kernel to measure memory latency values for a set of random strides. This micro-benchmark kernel runs in a pointer-chasing style, i.e., fetches a pixel and uses its value as 2D strides for subsequent access.

Step III: Cross-block stride feature extraction. In a 2D texture cache, data is typically organized in 2D blocks enabling data locality along both the width and height dimensions [15]. Due to this blocked data organization, data elements that look continuous from the logical (image) view may be stored far from each other in the physical memory if they belong to different data blocks. We can assume that accessing elements across blocks results in longer latency than accessing data within each block. Based on this assumption, we can build a machine-learning model to predict the relationship between data access patterns and access latency. We introduce the term **cross-block stride** for this machine learning model. It is defined as a function of the shape and size of the data block, and is *the stride that goes across distinct data blocks (under the assumed shape and size of the block)*. Our benchmarking collects the cross-block strides for various assumed data block shapes/sizes and the execution latency for each run. This information is used as the input feature of our subsequent machine-learning model.

Step IV: ML regression model fitting and analysis. Steps I-III are repeated multiple times to construct the training data (H, L), where H is the histogram of cross-block strides for assumed data block shapes/sizes and L is the profiled latency for each run. The collected training data are fed into an ML regression model based on the least squares method. For a given computation kernel and its associated data access pattern, we can first calculate the histogram of cross-stride accesses, and then use the model to predict the access latency.

3.2 Extension to Parallel Execution

As stated previously, multiple threads share the 2D cache. This section extends the previous benchmarking to a set of threads at both the warp and work-group levels.

3.2.1 Intra-warp spatial locality. This section studies the impact of threads within each warp on spatial data locality.

Empirical study settings. This experiment uses a single warp with varying numbers of threads – ranging from 1 to the maximum number supported (e.g., 64 on the mobile Adreno GPU). Each thread reads different data sections in a streaming manner. The entire data footprint is larger than the last level (L2) cache to rule out the impact of temporal data locality. This experiment is conducted on 4 kernels with representative data access patterns (as shown in Fig. 3). From the cache line reuse perspective, these kernels represent 4 cases with increasing reuse distances. Specifically, they are:

- ① Column major coalesced access: each thread reads data in column-major fashion and the threads collectively consume a full cache line immediately (cache line reuse distance of each thread is treated as 0),
- ② Row major: each thread reads each row or each cache line sequentially (cache line reuse distance being 1),
- ③ Block-2 major: each thread reads data in a 2-row block manner (cache line reuse distance being 2), and
- ④ Block-4 major: each thread reads data in a 4-row

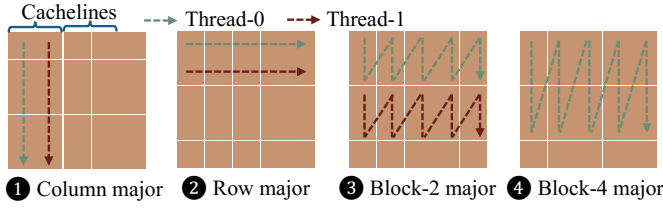


Figure 3: 4 different access patterns: with increasing cache line reuse distance from 0 to 4. In this example, each yellow block is a pixel, and two horizontal pixels occupy a cache line. Each thread consumes one pixel at a time.

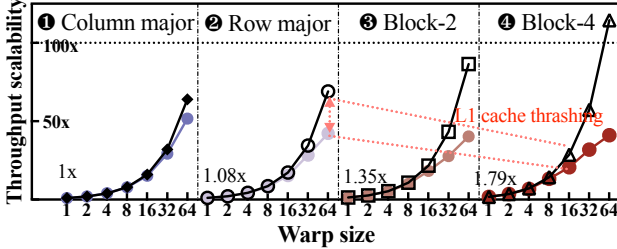


Figure 4: Scalability of each access pattern: restricted by reuse-distance. Black Line: perfect scalability for comparison. All results are normalized by 1 thread of Column major. For 1 thread, Row major, Block-2, and Block-4 outperform Column major by 1.08 $\times$, 1.35 $\times$, and 1.79 $\times$, respectively.

block manner (cache line reuse distance being 4). Each kernel runs 10 times and average memory throughput is measured. **Results Analysis.** Figure 4 shows the benchmarking results, where the x-axis corresponds to the number of threads in the warp, and the y-axis shows the measured memory throughput compared with *perfect scalability* (shown by black lines, calculated by $1_thread_throughput \times \#threads$). These results unveil two key findings: 1) With respect to single-thread throughput, Block access patterns (e.g., ③ and ④) achieve better performance because they are more consistent with the 2D texture memory block representation, i.e., result in fewer cross-block strides. 2) Regarding scalability, the performance of Block access patterns (e.g., ③ and ④) deteriorates quicker as the thread count increases. This is because Block results in larger cache line reuse distances, requiring more simultaneously active cache lines within a specific time window. It increases memory bandwidth utilization with a small number of threads as there are more concurrent memory requests. However, this leads to severe cache thrashing and contention when the number of threads increases.

We can observe that Block access patterns (e.g., ③ and ④) are friendly for low warp size scenarios; data access patterns with smaller cache line reuse distances (e.g., ① and ②) have the potential to achieve better performance as the thread count grows. A likely consequence is that for larger input shapes – where it is more likely that more threads

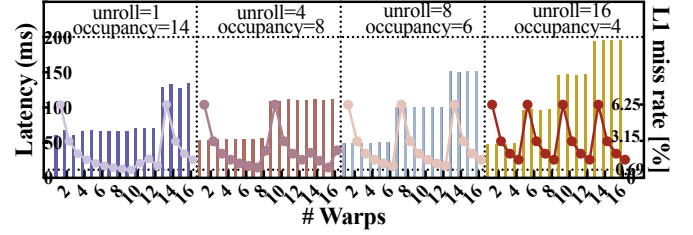


Figure 5: Intra-work-group temporal locality: Lines represent L1 miss rate using the right y-axis; Bars represent execution latency using the left y-axis.

are deployed – performance will be better with data access patterns that scale to a larger number of threads (e.g., ① and ②). Conversely, with smaller input shapes, where the code is likely to use few threads, an access pattern with higher single-thread performance (e.g., ③ and ④) is preferred.

Next, transitioning to a coarser-grained perspective, we study temporal data locality among threads (at intra-warp, inter-warp/intra-work-group, and inter-work-group levels). Due to limited space, we only describe experiments related to intra-work-group studies. Others are captured in our aforementioned online Supplementary Material (Sec. II).

3.2.2 Intra-work-group temporal locality and occupancy. This study aims to confirm that temporal data locality among warps within each work group can also help improve performance. Additionally, by analyzing execution latency results, this study identifies critical factors for mobile GPU occupancy modeling with varying work-group sizes.

Empirical study settings. This experiment uses a single work group, varying its size from 1 warp (64 threads) to 16 warps (1024 threads). All threads read the same data in a row-major pattern, exceeding the L2 cache capacity in all cases. An expensive function (e.g., log) is applied to the data read, with loop unrolling used to vary register pressure. Increased register usage reduces the number of concurrently executed warps, known as occupancy [64]. Unroll factors from 1 to 16 are tested, with occupancy ranging from 14 to 4. This experiment was repeated with other data access patterns, but each yielded similar conclusions.

Results Analysis. Fig. 5 shows the experimental results, where x-axis shows the work-group size in warps with all 4 different unrolling factors, the right y-axis shows the L1 texture cache miss rate, and the left y-axis shows the execution latency. For each unrolling case, the L1 miss rate decreases linearly when the work-group size is smaller than the occupancy, as all threads read the same data. When the work-group size exceeds the occupancy limit, the L1 miss rate exhibits an intriguing pattern: it initially increases, then drops to its lowest point, forming a recurring wave-like curve. We define *occupancy group* as a set of active warps that collectively fit within the available hardware resources (e.g., registers, and

execution units) of a GPU’s streaming multiprocessor (SM). This *chronic wave* behavior is observed for occupancy group. This indicates that warps within the same work group but in different occupancy groups cannot share data in the L1 cache. Additionally, even though the L1 cache miss rates are the same for different numbers of work groups, the latency values increase linearly with occupancy, resulting in a stepped latency curve. This leads to a conclusion that *warps in the same occupancy group execute concurrently, while those in different groups operate serially*.

Moreover, comparing execution latency among all unrolling cases, particularly between $unroll = 1$ and $unroll = 16$, with identical numbers of thread warps, reveals a trade-off between single-warp performance and occupancy. Both factors together influence overall performance. While unrolling uses more registers and improves performance per thread, it can reduce overall kernel performance by decreasing occupancy and increasing context switching overhead as the work-group size increases.

We now summarize the various observations. First, intra-work-group temporal locality exists, offering more opportunities for using Block access patterns (e.g., ③ and ④). Second, kernels with different input/output shapes/sizes may prefer different sizes of work groups due to occupancy considerations. For smaller data sizes, kernels favor smaller work groups with unrolling for improved single-thread performance. For larger ones, kernels prefer larger work groups without unrolling to maximize occupancy.

4 Performance model

Based on the benchmarking studies of the mobile GPU architecture introduced above, we developed an analytical model for performance prediction. The goal of this model is to enable (memory-related) performance optimizations and/or selection of better versions of code.

Fig. 6 summarizes the basic idea of this model and Table 1 summarizes all its parameters. The static information input to the model, which is independent of the program being executed, includes cache capacity (number of lines C) and the number of streaming processors (SPs). Parameters specific to the kernel but independent of the implementation include operator types (OP) and output shapes (O). The parameters that can vary with implementation include the data access pattern ($\mathcal{A}$), the work-group size $\mathcal{G} = b_x \times b_y \times b_z$ (x, y, z are three dimensions of the work group), and the thread-level tiling size ($\mathcal{T}$) which is equivalent to the unrolling factors mentioned in Section 3.2⁵. Specifically, each warp ($\mathcal{W}$) contains a fixed number of threads (64 on Adreno GPUs), with associated work-group size and shape $\mathcal{G}$. This performance model

⁵Our implementations empirically unroll all the tiling loops, hence tiling size equals the unrolling size.

Table 1: TModel parameters. Var.: variable, Con.: constant.

Var.	$\mathcal{A}, H_s$	Access pattern, cross-block strides histogram, reuse-distance
	$\mathcal{G}, \mathcal{W}$	Work-group size, warp shape
	$\mathcal{T}$	Tiling/unrolling
Con.	OP, O	Operator type, output shape
	SPs, R	#Shader cores, #Registers per core
	C, β	#Cache lines, LR modeled texture represent

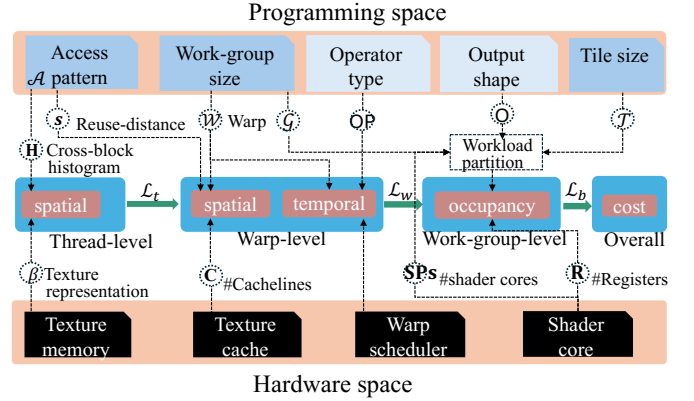


Figure 6: Performance model overview. The model is built from left to right by taking corresponding parameters.

enumerates all possible combinations of implementation-dependent factors, calculates the execution cost (cost) for each of them, and chooses the combination with the minimum cost. As we will describe in the next section, this performance model, though broader in its applicability, has already been integrated with a compiler that generates code parameterized with these parameters. Thus, by choosing parameters, the model enables optimized code for each kernel.

More specifically, the performance model consists of four components: 1) at the *thread level*, it models the texture memory latency for varied access patterns – it utilizes the results of the cross-block stride study from Section 3.1; 2) at the *warp level*, it models how texture memory latency changes for different access patterns as threads count grows – this component utilizes the scalability study from Section 3.2.1; 3) at the *work-group level*, it models the impact of occupancy and thread workload partitioning on parallelism within each work group and across work groups based on the study from Section 3.2.2; and 4) the *final level*, where the data access latency (first 2 components) and parallelism model (third level) are integrated to predict the execution cost of a given kernel. Please note: computation-related aspects (e.g., arithmetic intensity) are not considered in this model since our goal is to choose between different implementations of one operation – these implementations are expected to be identical with respect to these parameters.

Thread-level. At the thread-level, the cross-block stride information obtained earlier is summarized as a histogram

H capturing the frequency of strides across various block sizes. Our performance model takes as input the histogram and applies the fitted linear regression model introduced in Section 3.1 to predict texture memory access latency $\mathcal{L}_t(\mathbf{H})$. Specifically, the machine learning model described in the previous section results in a vector β of the same length, and each value is a weight that is to be multiplied by the corresponding element in the access pattern vector $\mathbf{H}$ (i.e. a dot product is performed). Thus, the predicted thread-level latency is given by:

$$\mathcal{L}_t(\mathbf{H}) = \langle \beta, \mathbf{H} \rangle. \quad (1)$$

Warp-level. The warp-level step maps the thread-level data access latency $\mathcal{L}_t$ computed above to the warp-level data access latency $\mathcal{L}_w$. This calculation should consider different *warp shapes*, marked as $\mathcal{W} = (w_x, w_y)$, and *warp size*, marked as $|\mathcal{W}| = \min(w_x \times w_y, 64)$. The w_x and w_y are configurable and subject to the work group size [61]. A detailed description of warp shape and thread mapping is in our on-line Supplementary Material (Sec. III). If there is no resource contention, $\mathcal{L}_w = \mathcal{L}_t$. However, as discussed in Section 3.2.1, the warp-level performance is restricted by two key factors: the texture cache capacity (i.e., the number of cache lines, denoted as C), and intra-thread reuse distance (s). The impact of the latter is dependent on the data access pattern (as shown in Fig. 4). When $s \times |\mathcal{W}| > C$, cache thrashing/contention occurs, thus the memory latency will increase. The effect of cache thrashing on increased latency is denoted by D . Then, the predicted texture memory access latency is:

$$\mathcal{L}_w(\mathcal{W}, \mathbf{H}, s) = \mathcal{L}_t(\mathbf{H}) \times D^{\lceil \frac{(|\mathcal{W}| \times s - C)}{C} \rceil}. \quad (2)$$

Intuitively, if $|\mathcal{W}| \times s \leq C$, $\lceil \frac{(|\mathcal{W}| \times s - C)}{C} \rceil$ equals to zero and cache thrashing decay rate is set to one (i.e. there is no cache thrashing). However, when the value is higher, the exponential term captures the slowdown observed. In further explaining this, if we take log on both sides, we get $\log \mathcal{L}_w - \log \mathcal{L}_t = \lceil \frac{(|\mathcal{W}| \times s - C)}{C} \rceil \log D$. The expression is derived by collecting data access latency values with different warp shape and applying the least squares method to solve for $\log D$. The value of D remains constant for each mobile GPU and does not vary per application, ensuring stability across different workloads.

For operators with a single input (e.g., SoftMax) the above formula is sufficient to calculate the warp-level texture memory access latency⁶. For operators with more than one input (e.g., MatMul), the memory access latency for each input tensor needs to be accumulated, as the contention for cache may be across different input tensors. Assuming we have i inputs, the general formula is:

$$\mathcal{L}_w(\mathcal{W}, \mathbf{H}, s) = \sum_i \mathcal{L}_t(\mathbf{H}_i) \times D^{\lceil \frac{(\sum_i \mathcal{W}_i \times s_i - C)}{C} \rceil}. \quad (3)$$

Work-group-level. As discussed, the work-group-level component of our cost model considers two key factors affecting execution cost: *occupancy* and *workload partitioning*.

Occupancy refers to the maximum number of warps that can be concurrently executed by each stream processor [64], as also discussed in Section 3.2.2. For example, each work-group on an Adreno GPU logically consists of up to 1024 threads, which is equivalent to 16 warps. Each work-group is exclusively assigned to one shader core, where registers are allocated for each thread for private usage. Hence, occupancy may vary depending on the register usage of each thread and overall register resources. Exceeding the occupancy over what can ensure sufficient register resources results in serial execution. The occupancy can be calculated as:

$$\text{Occupancy} = \mathbf{R} / \text{Register Footprint}(\mathcal{T}), \quad (4)$$

where Register Footprint depends on the computation kernel and thread-level tiling size ($\mathcal{T}$), a factor that is independent of the input size. Elaborating on this, as also discussed in Section 3.2.2, *unrolling* (or tile size) affects the register usage of each thread. Increasing the tile size leads to higher register usage per thread, which improves individual thread performance but decreases occupancy. The register footprint can be profiled offline with negligible overhead [54].

Workload partition refers to partitioning of the entire computation to each thread (in most implementations it is based on the output shape). We calculate the number of work groups (WGs) as follows:

$$\text{WGs} = \prod_j \left\lceil \frac{\mathbf{O}_j}{\mathcal{T}_j \cdot \mathcal{G}_j} \right\rceil. \quad (5)$$

Here, j is the index of the output dimension, $\mathbf{O}_j$ is the corresponding dimension size, $\mathcal{T}_j$ and $\mathcal{G}_j$ are the thread-level tile size and work-group size on the j th dimension, respectively.

As demonstrated in Sec. 3.2.2, the concurrency capability at the warp level is determined by both the occupancy and the number of streaming processors. To model this, we calculate the total number of *occupancy group* across all streaming processors, denoted as $\mathcal{L}_g$:

$$\mathcal{L}_g(\mathcal{T}, \mathcal{G}) = \left\lceil \frac{\text{WGs}}{\text{SPs} \times \text{occupancy}} \right\rceil. \quad (6)$$

Integration. The overall performance model for a kernel is expressed as $\mathcal{L} = \mathcal{L}_w \cdot \mathcal{L}_g$, encompassing both warp-level and work-group-level factors. Specifically, we define it as:

$$\sum_i \mathcal{L}_t(\mathbf{H}_i) \times D^{\lceil \frac{(\sum_i \mathcal{W}_i \times s_i - C)}{C} \rceil} \cdot \left\lceil \frac{\prod_j \lceil \frac{\mathbf{O}_j}{\mathcal{T}_j \cdot \mathcal{G}_j} \rceil}{\text{SPs} \times \text{occupancy}} \right\rceil. \quad (7)$$

⁶2D texture memory is read-only, so output write latency is not considered.

Applicability and extensibility. TMMoel is broadly applicable across diverse scenarios. Given its lightweight design and analytical modeling approach, we discuss its potential extensions to two other cases: dynamic shapes and irregular workloads.

Dynamic shapes. TMMoel explicitly accounts for input shapes, enabling it to handle dynamic input variables effectively. Owing to its analytical formulation and low overhead, TMMoel offers a greater advantage over existing methods such as auto-tuning (e.g., Table 5). It can adapt to dynamic input shapes by efficiently recomputing predictions at run-time.

Irregular workloads. Our current work focuses on optimizing dense and regular kernels, which are predominant in modern DNNs. Extending precise occupancy modeling to irregular workloads is part of our future work. This is a fundamentally challenging task due to data-dependent access patterns (e.g., sparse matrices) and additional complexity from unpredictable thread block scheduling. Notably, vendor documentation [55] states: “For Adreno GPUs, developers cannot assume the launch order of workgroups or warps in SPs.”

5 Model-based Code Optimization

We implement the cost model as a performance modeling module and integrate it with a state-of-the-art DNN compilation and acceleration framework (DNNFusion [47]) for mobile. Similar to other DNN compilation frameworks [9, 21], DNNFusion leverages a tensor-based DSL [48] and compute/schedule separation principle as Halide [57] to generate multi-versions of kernel codes in OpenCL. Halide [57] is a domain-specific language for image and tensor computations whose core design separates algorithm specification (what to compute) from scheduling (how to compute). We follow a similar design principle, allowing developers to explore diverse code transformations—such as loop tiling, unrolling, and memory layout changes—without modifying the underlying computation logic. To support data access patterns previously discussed, we extend DNNFusion by adding necessary schedules, data storage, and assisted code templates. We also extend DNNFusion to support backward propagation, thus supporting on-device DNN training.

The cost model can help identify the optimal combination of *data access pattern* (and its corresponding memory allocation), *output tile size* (or loop unrolling factor), and *thread configuration* (including warp shape and work-group shape). The first two are *implementation factors* and will be used by our compiler to generate the corresponding schedule and kernel code. The third is a *configuration factor* and will be passed to the generated kernel code as a parameter.

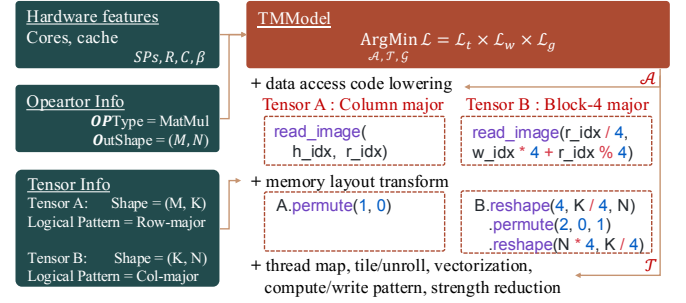


Figure 7: Compiler workflow. A MatMul example that generates two data access patterns: Column and Block-4 major. Both data access pattern and memory layout transformation are needed. Other optimizations happen subsequently.

```

1 __constant sampler_t sampler =
2   CLK_NORMALIZED_COORDS_FALSE | CLK_ADDRESS_CLAMP |
3   CLK_FILTER_NEAREST;
4 __kernel void matmul(
5   __read_only image2d_t A,
6   __read_only image2d_t B,
7   __write_only image2d_t C) {
8   // thread mapping
9   const int w_idx = get_global_id(0);
10  const int h_idx = get_global_id(1);
11
12  float4 c = 0;
13  // REDUCE_SIZE is determined by the tile/unroll schedule
14  for(short r_idx=0; r_idx<{{REDUCE_SIZE}}; r_idx++) {
15    float4 a = read_imagef(
16      A, sampler, (int2)(h_idx, r_idx));
17    float4 b = read_imagef(
18      B, sampler, (int2)(r_idx/4, w_idx*4+r_idx%4));
19
20    // Compute pattern is determined by vectorize schedule
21    c = mad(a.x, b, c);
22    c = mad(a.y, b, c);
23    c = mad(a.w, b, c);
24    c = mad(a.z, b, c);
25  }
26  // OUT_IDX is determined by the write pattern
27  // schedule, which defines the output tensor layout
28  // optimized for downstream consumers.
29  write_imagef(
30    C, (int2)({{OUT_IDX(w_idx, h_idx)}}), c);
31 }

```

Listing 1: MatMul Kernel Example. This kernel is generated after applying the data access pattern schedule. Highlighted placeholders (e.g., REDUCE_SIZE, OUT_IDX) are determined by subsequent schedules such as tiling, unrolling, and write pattern.

It is worth noting that although we have deployed TMMoel for code generation and optimization in DNNFusion, it equally applies to both other DNN compilation frameworks like TVM [9] for code generation and library-based DNN acceleration frameworks like MNN [30] for kernel selection on mobile GPUs with 2D memory.

Figure 7 shows an indicative workflow of using TMMoel to optimize MatMul. This example also shows more details

about the schedule and code lowering for several data access patterns. These schedules are also supported by Halide [57], specifically involving the design of appropriate *domain orders* (and loop transformations) described in Section 3.2 of [57]. The correctness of these schedules is guaranteed because the change in data access pattern and data storage order still maintains the original logical data access order. The schedule describing the data access pattern is decided before all other schedules (e.g., *tiling* and *unrolling*) and supports individual kernel and fused kernel in DNNFusion [47]. Listing 1 presents the corresponding lowered OpenCL code, illustrating the impact of different access patterns on the generated kernel. This example omits subsequent schedules (e.g., *tiling*) as they are natively supported by Halide.

6 Evaluation

This section assesses the efficacy of TMMoel by comparing it with *four* state-of-the-art mainstream product-level mobile DNN acceleration frameworks (and their associated performance modeling/tuning methods): MNN [30], CLML [1], TFLite [2], and TVM [9]. We also compare against a more general-purpose GPU performance modeling method (PPT-GPU [3]). Both individual kernels and end-to-end on-device DNN training are used as target workloads. More specifically, the *evaluation objectives* are as follows: 1) demonstrating that TMMoel can lead to better kernel performance, outperforming all baseline methods (Section 6.2); 2) showing the efficacy of TMMoel in end-to-end full on-device DNN training (Section 6.3); 3) performing an *ablation study*, i.e., understanding the effectiveness of different components of our cost model, including verifying the model’s prediction accuracy at each level (Section 6.4); 4) verifying the cost of TMMoel-based code optimization by comparing with other baselines, validating the portability and generality of TMMoel by showing its performance on different mobile GPUs, and evaluating its inference speed against other inference frameworks (Section 6.5).

6.1 Experiment Settings

Baselines. Table 2 summarizes the key aspects of five baselines used in our evaluation. Particularly, MNN, CLML, and TFLite explore different work-group sizes using brute-force search and select the best one. The work-group size (G) configuration is independent of kernel implementations, but it is important for achieving optimal performance. TVM’s auto-tuning approach predicts kernel latency by an ML-based cost model and searches optimization factors (e.g., tile sizes) by generating code and profiling. TFLite performs heuristic kernel selection based on input shapes. TMMoel is also compared with a general GPGPU memory modeling approach (originally for NVIDIA GPUs), PPT-GPU [3] that calculates

Table 2: Baselines characterization. A: data access pattern; T: thread tile; W: Warp shape; G: Work-group size.

Framework	MNN	CLML	TVM	TFLITE	PPT-GPU	TMMoel
Perf. model	searching	searching	ML-Based	searching	analytical	analytical
Opt. target	G	G	T, G	T, G	A	A, T, W, G
m-GPU train	Yes	Yes	No	No	No	Yes

cache miss rates given the memory trace/data access order. Cache miss rate is directly related to the warp shape (temporal locality), so PPT-GPU implicitly searches warp shape. Compared to TMMoel, PPT-GPU does not target 2D texture memory, it requires extensive knowledge of GPU hardware parameters (e.g., the cache associativity), and relies on memory trace to calculate reuse distance. On a mobile GPU hardware parameters are not documented, and no tracing tools are available to extract address information. To overcome this, we follow the process outlined in earlier work [62] to generate memory traces, in addition to using the information on associativity collected in our experiments [59, 70].

Workloads. TMMoel is evaluated on representative individual DNN operators and end-to-end DNN on-device training: *Individual kernels/operators*. The evaluated kernels include Matrix Multiplication (MatMul), 2D Convolution with strides of 1 (Conv2Ds1) and 2 (Conv2Ds2), 2D DeConvolution with strides of 1 (DeConv2Ds1) and 2 (DeConv2Ds2), 3D Convolution (Conv3D), 2D Depthwidth Convolution (DwConv2D), SoftMax, and LayerNorm. Their typical application DNNs and ranges of input sizes for our evaluation are summarized in the left half of Table 3. To optimize these kernels, TMMoel explores the best access patterns among five pre-defined ones (as shown in Fig. 3) independently for each input: Column-major, Row-major, and three Block-major ones with reuse distance of 2, 4, and 8. Furthermore, varying tiling sizes are considered for each pattern.

On-device training. Training of four representative DNN models is also used for TMMoel’s evaluation: we use three classic Convolutional Neural Networks with different major kernels, i.e., MobileNet [60] with DwConv2D, ResNet-18 [23] with Conv2D, and Fsrcnn [16] with DeConv2D, and a popular vision transformer Vit [17] with MatMul. We use on-device training for our *full application* evaluation because: 1) it becomes increasingly important for up/downstream applications such as federated learning [42]; 2) it is more challenging than inference – specifically, the forward-backward computational graph has more shape variations and more complex data dependencies; 3) its execution on mobile GPUs remains an open challenge [13, 73]. Our evaluation focuses on the kernel optimization aspects of on-device training without resorting to algorithm optimizations (e.g. the ones that may sacrifice numerical precision [36] or memory space [37]) – these optimizations are orthogonal to TMMoel.

Table 3: Individual kernel performance improvements over baselines. Numbers outside and within parentheses denote the *average* and *peak* (tagged w/ $\uparrow$) speedup over baseline, respectively. '-' means this framework does not support this kernel on mobile GPU. The third column specifies the input size range of each input: MatMul has two inputs, each with each dimension size ranging 2^4 - 2^{12} (double every time). For Conv, two sets are for input matrix and kernel size of weight matrix, respectively.

Kernels	Typical applications	Shape range	Avg. speedup over				
			MNN	CLML	TVM	TFLite	PPT-GPU
MatMul	Transformers	2^4 - 2^{12}	2.57 (3.73 $\uparrow$)	6.04 (9.17 $\uparrow$)	5.34 (11.9 $\uparrow$)	1.20 (3.22 $\uparrow$)	1.47 (2.78 $\uparrow$)
Conv2Ds1	CNNs	2^{4-9} , 3 – 7	1.35 (1.57 $\uparrow$)	2.06 (2.55 $\uparrow$)	1.18 (1.80 $\uparrow$)	1.12 (1.53 $\uparrow$)	1.07 (1.38 $\uparrow$)
Conv2Ds2	Downsample	2^{4-9} , 3 – 7	1.19 (1.82 $\uparrow$)	1.42 (2.49 $\uparrow$)	1.42 (4.36 $\uparrow$)	1.20 (1.63 $\uparrow$)	1.16 (1.81 $\uparrow$)
DeConv2Ds1	Conv backward	2^{4-9} , 3 – 7	2.16 (2.41 $\uparrow$)	2.32 (2.70 $\uparrow$)	1.71 (4.13 $\uparrow$)	1.45 (3.49 $\uparrow$)	1.79 (2.41 $\uparrow$)
DeConv2Ds2	Upsample	2^{4-9} , 3 – 7	2.18 (2.42 $\uparrow$)	6.57 (7.45 $\uparrow$)	3.96 (8.60 $\uparrow$)	1.20 (2.53 $\uparrow$)	1.27 (2.42 $\uparrow$)
Conv3D	Action recognition	2^{4-9} , 3 – 7	–	–	6.62 (8.67 $\uparrow$)	–	4.47 (8.67 $\uparrow$)
DwConv2D	CNN	2^{4-9} , 3 – 7	2.49 (3.41 $\uparrow$)	1.62 (4.11 $\uparrow$)	2.10 (4.85 $\uparrow$)	1.76 (3.24 $\uparrow$)	2.13 (3.41 $\uparrow$)
SoftMax	Transformers	2^4 - 2^{12}	4.85 (8.88 $\uparrow$)	3.62 (12.5 $\uparrow$)	2.32 (4.36 $\uparrow$)	1.31 (8.83 $\uparrow$)	2.28 (8.88 $\uparrow$)
LayerNorm	Transformers	2^4 - 2^{12}	3.43 (4.34 $\uparrow$)	–	7.77 (18.2 $\uparrow$)	3.64 (14.1 $\uparrow$)	2.53 (4.34 $\uparrow$)
Geometric Mean			2.31 (3.08 $\uparrow$)	2.86 (4.82 $\uparrow$)	2.91 (6.06 $\uparrow$)	1.48 (3.61 $\uparrow$)	1.83 (3.30 $\uparrow$)

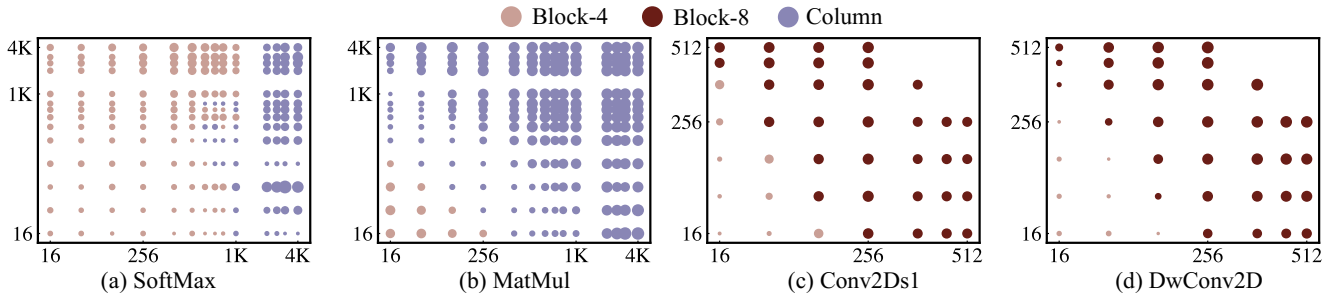


Figure 8: Output shape impacts kernel selection. Circle colors represent different implementations. Circle sizes represent the speed improvements over the base implementation which is a vanilla version without tiling or data access pattern optimizations, but with fine-tuned work-group size. The x- and y-axis denote the height and width of output geometry.

Evaluation environment. Evaluations are carried out on the Qualcomm Snapdragon 8 Gen 2 SoC (SND Gen2) [56] with a Qualcomm Adreno 740 GPU, and version 3.0 of the OpenCL driver. Additionally, MediaTek D1100 [43] SoC with Mali G77 GPU is used to demonstrate portability. For individual DNN kernels, each experiment was run 10 times, with the ‘warm-up’ run excluded, and the average execution time of other runs is reported. The confidence interval is omitted for readability, as the variations were low. For the on-device training experiments, a consistent configuration is applied for each framework and test model: the batch size is set to one, and Stochastic Gradient Descent (SGD) is used to update model parameters once every 32 batches. This approach is commonly used [58, 68] to save memory space while maintaining the same training effectiveness as a standard batch size of 32. The reported latency is the average execution time for one batch, calculated over 32 batches. To maintain training accuracy, floating-point 32 is used for on-device training evaluation (as well as for individual kernel evaluation).

6.2 Overall Performance

Table 3 summarizes the generated code speedup enabled by TMMoel over all five baselines. The numbers outside and inside parentheses denote the average and peak speedup respectively of TMMoel over the corresponding baseline across all input sizes for each kernel. For example, for MatMul, TMMoel achieves an average speedup of 2.57 $\times$ and a peak speedup of 3.73 $\times$ over MNN across all input sizes (size of each dimension ranges from 2^4 to 2^{12}). For all individual kernels, the geometric means of the average speedsups of TMMoel over MNN, CLML, TVM, TFLite, and PPT-GPU are 2.31 $\times$, 2.86 $\times$, 2.91 $\times$, 1.48 $\times$, and 1.83 $\times$, respectively, and the geometric means of the peak speedsups are 3.08 $\times$, 4.82 $\times$, 6.06 $\times$, 3.61 $\times$, and 3.30 $\times$, respectively, demonstrating TMMoel’s efficacy.

TMMoel outperforms other frameworks for two main reasons: First, it systematically considers multiple optimization factors (data access patterns, thread-level tiling, and warp and work-group level thread mapping) simultaneously, while others mainly focus on one or two of them (see Table 2). For

Table 4: On-device training performance comparison.

Models	Major Kernels	Latency (ms)		
		MNN	DNNF	TMMoel
MobileNetV2 [60]	DwConv2D	1916	41	29
ResNet-18 [23]	Conv2D	88	66	48
Fsrrnn [16]	DeConv2D	921	57	32
ViT [17]	MatMul	2105	1224	575

example, MNN only supports work-group fine-tuning without exploring kernel implementation options. Second, and more importantly, TMMoel takes into account the 2D texture memory (and cache) and couples it with thread mapping and occupancy considerations. Examining results from other frameworks – although PPT-GPU models cache behavior, it does not take texture representation, occupancy, or output shape into account for memory access and thread mapping optimizations. TFLite selects different implementations for specific input/output shapes, so it performs the best among all baselines. However, it also misses to consider texture representation and occupancy in a nuanced manner. CLML searches only a limited set of candidates and misses many optimization opportunities. Finally, TVM lacks specialized templates for many kernels on mobile GPUs and fallbacks to using a generic template.

Output shapes impact kernel configuration selection.

An advantage of TMMoel is the consideration of output shape. Fig. 8 verifies this by showing the distribution of best-performing implementations for four representative kernels - SoftMax, MatMul, Conv2Ds1, and DwConv2D⁷. For example, for SoftMax and MatMul, with small input shapes and low parallelism, Block-major access is preferable to achieve low single-thread data access latency ($\mathcal{L}_t$). However, as the input size increases, using Block-major (with large s) and more threads (larger W) would cause cache thrashing, resulting in degraded performance (increased $\mathcal{L}_w$ thus $\mathcal{L}_g$). For DwConv2D and Conv2D, Block-major access still results in ideal performance as their input sizes grow because their input of the *Weight* matrix has high data reuse (and temporal locality), effectively reducing cache thrashing as thread count grows. For readability, Fig. 8 only shows the selection of data access patterns, omitting the tiling factor determination.

6.3 End-to-End DNN On-Device Training

Among all five baselines, MNN is the state-of-the-art framework to support end-to-end DNN on-device training on mobile GPUs. Table 4 compares TMMoel with MNN on four representative DNN models with varied major kernels. TMMoel achieves 1.83 – 66.1 $\times$ speedup over MNN for two main reasons: First, TMMoel is built on DNNFusion [47] (with the latest

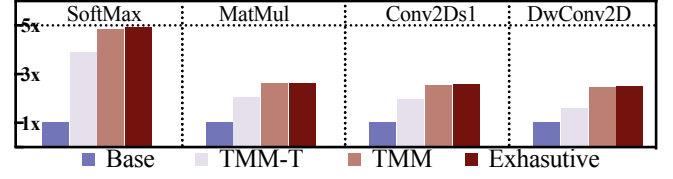


Figure 9: Performance breakdown with varied model outputs. TMM refers to TMMoel and TMM-T uses predicted tiling factor only for code opt. Base is a vanilla version as Fig 8 while Exhaustive refers to a fully optimized version with exhaustive search. y-axis is the speedup over Base.

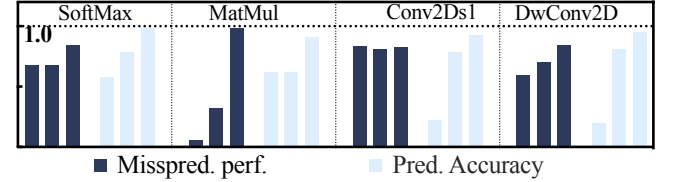


Figure 10: Prediction accuracy and average latency for all miss-prediction cases with each level of TMMoel for four kernels. The bars of each group from left to right represent thread-/warp-/integrated-level of TMMoel.

memory optimization [49]), which has more advanced operator fusion and memory operation elimination than MNN. Second, TMMoel further optimizes DNNFusion with the newly designed 2D texture-aware analytic model to generate more efficient operators (or kernels), with a nuanced selection of data access patterns, tiling, and other GPU configuration parameters. Table 4 also compares TMMoel with DNNFusion, showing 1.38 – 2.53 $\times$ speedup on these four DNN models.

6.4 Performance Understanding

To further understand TMMoel’s performance, this section reports an ablation study on the four representative kernels we previously used for the detailed evaluation.

Breakdown of performance improvement with varied model outputs. Fig. 9 shows a performance breakdown with varied model outputs: Base is a vanilla version w/o tiling or data access pattern optimizations, but w/ fine-tuned work-group size. TMM-T uses predicted tiling sizes from TMMoel only for code optimizations. TMM uses both predicted tiling and data access patterns from TMMoel. Exhaustive shows the upper bound achieved by exhaustively exploring the search space (A, T, W, G). This study shows that both the access pattern and tiling are essential factors for performance tuning on mobile GPUs and TMMoel can achieve performance close (within around 1%) to that of exhaustive search. **Model prediction accuracy study.** We focus on 1) the prediction accuracy of TMMoel, and 2) what happens when TMMoel produces inaccurate configurations. Fig. 10 summarizes the prediction accuracy of each level of TMMoel

⁷Other kernels show similar trends.

Table 5: Tuning latency comparison. The tuning latency is the total time for finding the best configurations for Conv2Ds1 with all input sizes in Table 3.

Framework	MNN	CLML	TVM	TFLite	PPT-GPU	TMMoel
Tuning Lat. (s)	73.4	10.4	234	552	771	1.92

for these four kernels⁸, and the average latency of the miss-predicted cases by normalizing it to the latency of the best configuration with exhaustive search. These prediction accuracy results indicate that as more and more factors in TMMoel are considered, its prediction accuracy improves, achieving 94.1% accuracy on average for these four operators when the entire performance model is used, demonstrating TMMoel’s efficacy, and the necessity of each of its components. The miss-prediction performance demonstrates that for certain cases, even if TMMoel produces a wrong prediction, the predicted configuration achieves performance similar (within 83.8% – 98.4%) to the correct one.

6.5 Other Experiments

Tuning Latency. Table 5 compares the total tuning latency of TMMoel and all five baselines. We identify the best configurations for Conv2Ds1 with all input sizes (shown earlier in Table 3). Compared with all baselines, TMMoel accelerates the tuning stage by 5.42 – 402×. MNN, CLML, and TFLite are slower because of the exhaustive search and execution of the code candidates. TVM’s latency is also high because it iteratively profiles generated code on the mobile device and trains ML models. Although TMMoel also needs a benchmarking stage with a regression model (as described in Section 3), this is performed once for each architecture, rather than for each kernel (like TVM). PPT-GPU, even though being an analytical performance model, is slow because it calculates the reuse distance with an expensive analysis of memory traces [46].

Portability Study. Table 6 compares the on-device training latency of MNN and TMMoel on another GPU architecture, Mali G77, with one CNN (MobileNetV2) and one

Table 6: Portability study

Models	Latency (ms)	
	MNN	TMMoel
MobileNetV2	4060	59
ViT	5987	1936

Transformer (ViT). Other DNN models show similar trends. The Mali GPU family differs significantly from Adreno GPUs – it has a smaller warp size (32 threads), a larger texture size (8KB), more shader cores (9), a distinct texture representation, and features separate on-chip texture and data caches. With these critical differences, targeting this device demonstrates the portability

Table 7: Inference performance comparison. All frameworks use FP16 precision and run on a Snapdragon 8 Gen 2 platform with an Adreno 740 GPU.

Models	Latency (ms)		
	MNN	DNNF	TMMoel
ViT	533	277	88
ResNet-18	10.9	10.1	8.6

of TMMoel. To maximize the utilization of both data paths, we empirically employ a mixed-use approach where tensor inputs and outputs are handled with Image2D objects, while model weights are stored in Buffer objects. The evaluation results show that TMMoel outperforms MNN by 68.8× and 3.09× on MobileNetV2 and ViT, respectively.

Inference Latency. TMMoel is a general-purpose performance model applicable to both inference and training workloads. To evaluate its effectiveness for inference, Table 7 presents latency results on ViT and ResNet-18. TMMoel achieves significant improvements, outperforming MNN by 6.06× on ViT and 1.27× on ResNet-18, and surpassing DNNF by 3.15× and 1.17×, respectively.

7 Related Work

Texture memory and micro-benchmarks. Texture cache was originally developed for faster rendering on graphic processors [15, 22, 26]. There have been several micro-benchmarking studies to understand the performance of systems with such a cache. The work by Wong et al. [70] and Mei et al. [44] only used 1D data (Image1D) to dissect texture cache organization – generalizability to Image2D was left unclear. Romou [39] profiled cache hierarchy on mobile SoCs, however, the 2D spacial locality of texture memory was not studied. Our work distinguishes itself by studying texture representation and its impact on both spatial and temporal locality.

Cache and locality analysis for 1D memory. Reuse distance [14, 31] is a well-known metric to model cache behavior, and is studied based on an ordered memory access trace. Nugteren et al. [50] extends reuse distance to GPU cache sets by modeling GPU’s parallel execution paradigm and predicting cache miss rates. PPT-GPU-Mem [3], on the other hand, leverages a binary instrumentation tool NVBIT [63] to provide a reliable memory access trace to calculate reuse distance. Applying reuse distance to mobile platforms with texture cache is not straightforward, because no trace tools are available and the texture data representation is patented.

Cost models and tensor computing. Models of performance execution are designed for numerous objectives. Roofline model [69] identifies the bottleneck of hardware resources. Gables [24] retargets the Roofline model to model each accelerator on a SoC, apportions work concurrently

⁸A more detailed study is presented in our online Supplementary Material, Sec. IV.B, specifically, Fig.13 due to limited space.

among different accelerators, and calculates a SoC performance upper bound. Li et al. [38] extract features from the architecture and the program to train a generalizable cost model. MAESTRO [34] models DNN accelerators and estimates execution latency for a given the DNN model. Inspired by TVM [9] and Ansor [78], several works employ machine learning methods to predict inference latency and auto-tune DNN kernels [75, 77, 81]. TSM2 [8] designs a cost model to optimize tiling parameters for matrix multiplications with tall-and-skinny inputs. Babalad et al. [6] train machine learning models to select tile sizes and loop orders for CPUs. Jang et al. [29] optimize memory type selection by analyzing data access patterns in existing code. In contrast, our work focuses on generating kernels with efficient data access patterns tailored to the tensor shape and leveraging the 2D spatial locality of texture memory.

On-device training. There is a significant amount of work on DNN training on desktop GPUs [7, 11, 19, 74, 79], or other processors [27, 28, 72] – however, the challenges associated with distinctive features and resource limitations of mobile chips are not handled there. On the other hand, previous research on mobile devices has focused on optimizing on-device training performance through sparsification [76] and quantization [40, 73, 80]. Unlike these approaches, our methods do not compromise model accuracy. Additionally, a separate line of research primarily targets the orthogonal direction of memory consumption reduction [20, 40, 45, 67, 68].

8 Conclusion and Future Work

This paper presented a new performance modeling and optimization approach called TMMoel for mobile GPUs that takes 2D texture memory into account. TMMoel consists of three advances: 1) a micro-benchmarking and machine learning-based approach on top of a new concept (cross-block stride) to understand 2D spatial locality and analyze the impact of factors like warp size/shape, occupancy, and work-group size; 2) a comprehensive analytic model for choosing factors like data access pattern(s) and tile size(s) for a given kernel (and associated size/shapes), and 3) a complete system prototype that supports code optimization for both kernels and full DNN on-device training. TMMoel is extensively evaluated by comparing with four state-of-the-art mobile DNN frameworks and a general-purpose GPU performance model, achieving up to $3.61\times$ speedup on individual kernels and $66.1\times$ speedup for end-to-end DNN on-device training with as low as 0.25% tuning cost.

Based on TMMoel, our future work includes: 1) designing more advanced data layouts, and access patterns, 2) global layout selection, transformation, and memory elimination for DNN training, and 3) supporting irregular kernels (e.g., on sparse matrices or tensors).

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