

# A 95.6-TOPS/W Deep Learning Inference Accelerator With Per-Vector Scaled 4-bit Quantization in 5 nm

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**Abstract**—The energy efficiency of deep neural network (DNN) inference can be improved with custom accelerators. DNN inference accelerators often employ specialized hardware techniques to improve energy efficiency, but many of these techniques result in catastrophic accuracy loss on transformer-based DNNs, which have become ubiquitous for natural language processing (NLP) tasks. This article presents a DNN accelerator designed for efficient execution of transformers. The proposed accelerator implements per-vector scaled quantization (VSQ), which employs an independent scale factor for each 64-element vector to enable the use of 4-bit arithmetic with little accuracy loss and low energy overhead. Using a multilevel dataflow to maximize reuse, the 5-nm prototype achieves 95.6 tera-operations per second per Watt (TOPS/W) at 0.46 V on a 4-bit benchmarking layer with VSQ. At a nominal voltage of 0.67 V, the accelerator achieves 1734 inferences/s/W (38.7 TOPS/W) with only 0.7% accuracy loss on BERT-Base and 4714 inferences/s/W (38.6 TOPS/W) with 0.15% accuracy loss on ResNet-50 by using quantization-aware fine-tuning to recover accuracy, demonstrating a practical accelerator design for energy-efficient DNN inference.

**Index Terms**—Accuracy-efficiency trade-off, BERT, deep neural network (DNN) inference accelerator, quantization, transformers.

## I. INTRODUCTION

DEEP neural networks (DNNs) are now a foundational building block in a wide variety of application domains, including computer vision (CV) [1], natural language processing (NLP) [2], computer graphics [3], autonomous driving [4], and many others. While early DNNs generally employed convolutional neural network (CNN) architectures [5], the introduction of attention-based transformer architectures [6]

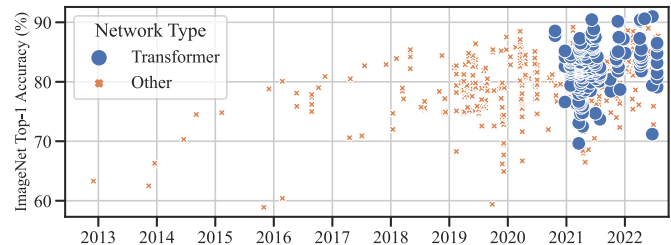


Fig. 1. Transformers achieve high accuracy on image classification [11].

and the subsequent success of very large transformer-based language models on natural language tasks [7] has motivated the development of transformers for CV as well [8] (see Fig. 1). These trends suggest that transformers will become an increasingly common workload across a wide range of applications and deployment targets.

The high arithmetic intensity and deterministic structure of DNN inference computation makes this workload an ideal candidate for dedicated hardware acceleration to improve energy efficiency [measured in tera-operations per second per Watt (TOPS/W)] and area efficiency (measured in TOPS/mm<sup>2</sup>) compared with software implementations on CPUs or GPUs. One common technique to improve hardware efficiency is quantization, in which data are represented as integers of eight or fewer bits, reducing both arithmetic and data movement costs. However, excessive quantization can introduce errors that impact the task accuracy of the DNN [9], and transformers are more sensitive than CNNs to quantization error [10].

This work presents a DNN inference accelerator that achieves very high energy efficiency while maintaining task accuracy on transformer workloads with a technique called per-vector scaled quantization (VSQ). After providing additional background in Section II, we describe our accelerator architecture and key design decisions in Section III. Section IV describes our approach to quantization-aware training (QAT) and mapping of transformers, key software workflows developed in tandem with the accelerator to maintain high accuracy. Section V highlights key measurement results, including an energy efficiency of 95.6 TOPS/W, a 5.8 $\times$  improvement over prior work.

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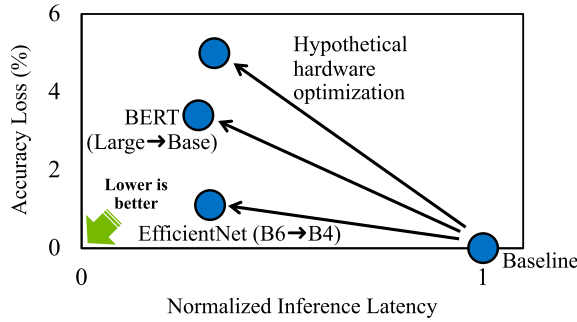


Fig. 2. Simpler network models can often outperform hardware optimizations. Latency data are gathered from [19] and [20].

## II. BACKGROUND

### A. Accuracy/Energy Trade-Off

Energy efficiency is a critical metric for DNN inference, both for power-constrained datacenter servers and battery-limited mobile devices. Prior DNN inference accelerators have proposed energy-saving techniques, such as quantization [12], sparsity [13], low-rank approximation [14], analog and in-memory computing [15], and approximate computing [16], all of which improve efficiency by either reducing the energy cost of each operation or reducing the number of operations required to complete the inference task. Each of these techniques introduces error in the computed result, and while DNN inference is resilient to small errors in computation [17], large enough errors can propagate to the network output, resulting in a reduction in accuracy. These approaches, therefore, introduce a trade-off between energy and accuracy, with more aggressive energy-saving measures resulting in larger errors and accuracy drops on a given network and task.

One challenge of realizing hardware techniques for energy efficiency improvement is that the amount of accuracy that can be sacrificed for improved efficiency is very small in practice. Consider a hypothetical hardware optimization that achieves a  $3\times$  reduction in inference latency while incurring a 5% accuracy loss. This may seem like a worthwhile trade-off, but as Fig. 2 shows, deep networks, such as BERT [2] and EfficientNet [18], can easily be downsized to shallower counterparts with a more favorable efficiency-accuracy trade-off, with no special hardware required. (We compare latency in this example, because benchmarking data are more readily available, but energy efficiency follows similar trends.) It is critical that the proposed energy-saving techniques have a negligible impact on task accuracy, so that they can outperform the straightforward energy-saving technique of employing a simpler network.

### B. Per-Vector Scaled Quantization

Reducing arithmetic precision by quantizing 32-bit floating-point (FP32) number representations to low-bitwidth integers is a powerful technique to save energy. The energy cost of data movement and storage scales linearly with bitwidth, while the cost of the multiply-and-accumulate (MAC) operations that comprise the core arithmetic of deep learning (DL)

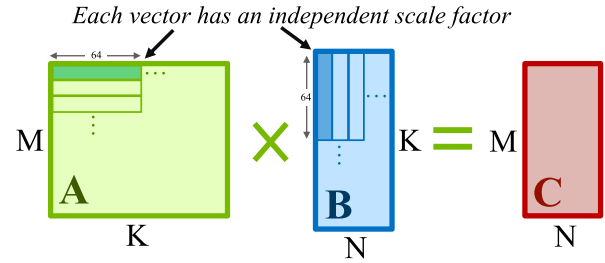


Fig. 3. MM with VSQ.

scales quadratically with input bitwidth. Typical quantization techniques employ a coarse scale factor granularity, using one common scale factor applied to many thousands of values within an input tensor. While per-tensor scaling is computationally inexpensive as the cost of the scale factor multiplication is amortized across many elements, it tends to require mapping a wider range of values to the low-precision representation, increasing quantization error. Quantization to 8-bit integers (INT8) can generally be realized with minimal ( $<0.1\%$ ) accuracy loss on most DNN architectures, but naive quantization to smaller representations can lead to catastrophic accuracy loss. For example, BERT-Large can be quantized to INT8 with only 1.5% accuracy loss using post-training quantization, but further reduction in precision to 4 bits (INT4) results in an unacceptable accuracy loss of 84% [21].

This work instead uses VSQ with fine-grained scale factors to enable high accuracy on transformers with 4-bit integer representations [21]. VSQ exploits the vector granularity of computation in many DNN inference accelerators. In addition to a coarse-grained per-output-column scale factor, VSQ quantizes data at a per-vector granularity within each input matrix in the matrix-matrix multiplication (MM), as shown in Fig. 3. The range that must be captured by each per-vector scale factor is much smaller than per-tensor scaling, greatly reducing quantization error (see Fig. 4). Per-vector scale factors require minimal hardware overhead, with one additional data word stored per input vector, and one additional scalar multiplication of the two vector scale factors for each vector dot product. The decreased quantization error from VSQ with 4-bit arithmetic (INT4-VSQ) enables minimal DNN accuracy loss on transformers compared with unquantized floating-point representations, while imposing only a small hardware overhead compared with INT4 math.

### C. Transformer Networks

Fig. 5 shows the structure of a typical transformer network, which consists of a sequence of transformer layers, each of which contains multihead attention and feed-forward blocks. The structure of transformers differs significantly from that of CNNs. Some MM operations are performed with pretrained weights that do not change during inference; these operations are similar to  $1 \times 1$  convolution (CONV) kernels. However, the batch matrix multiply (BMM) operations have no pretrained weights, with both inputs computed dynamically during runtime. While CNNs contain one softmax calculation, transformers perform softmax more frequently, once per

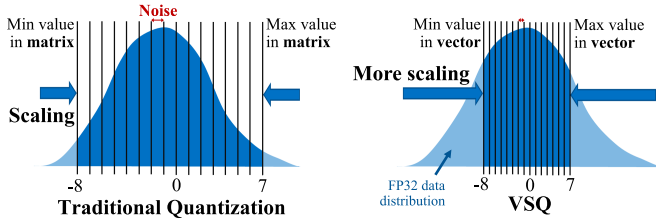


Fig. 4. VSQ reduces quantization noise.

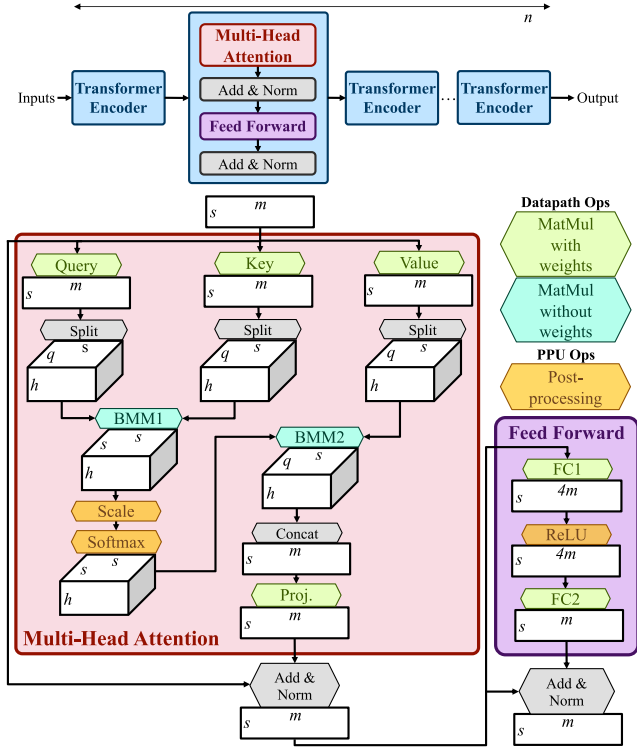


Fig. 5. Structure of transformer networks.

encoder layer. Dimensions of the transformer computations in BERT and DeiT [22] are listed in Table I (except for sequence length  $s$ , which varies with workload).

Note that we modify some of the computations within the standard transformer for better hardware efficiency. We replace the expensive nonlinear Gaussian error linear unit (GELU) operation with the simpler rectified linear unit (ReLU) function (the accuracy impact of this change is minimal after retraining). For simplicity, we also ignore some operations with low arithmetic intensity (shown in gray in Fig. 5) in our accelerator design and evaluation. In practice, these operations would be performed at a higher level of system hierarchy (see Section III-F).

Since MMs form a key building block for transformers, we use them as an example to illustrate different optimizations proposed in this work. For generality across MMs with and without weights, we denote the input matrices as  $A$  and  $B$  with dimensions  $M \times K$  and  $K \times N$ , respectively, with the output matrix  $C$  of dimensions  $M \times N$  (see Fig. 3).

#### D. Prior Work

Recent years have seen a proliferation of DNN inference accelerator designs. Digital accelerators exploit

TABLE I  
TRANSFORMER NETWORK PARAMETERS

Network	BERT-Large	BERT-Base	DeiT-Base	DeiT-Small
Model ( $m$ )	1024	768	768	384
Heads ( $h$ )	16	12	12	6
Query Size ( $q$ )	64	64	64	64
Layers ( $n$ )	24	12	12	12
MACs ( $\times 10^9$ )	123	35	8.5	1.1

quantization [12], [23], [24], [25], [26], [27], [28], sparsity [14], [23], [24], [27], and other energy-saving hardware optimizations [16], [29], while compute-in-memory accelerators make use of analog circuit properties to achieve very high reported energy efficiencies [15], [24], [25], [30], [31], [32]. However, many prior hardware accelerator proposals fail to evaluate task accuracy entirely [15], [27], [31], evaluate their proposals only on very small networks [25], [29], [30], or suffer large ( $>1\%$ ) drops in accuracy on realistic workloads [24], [32]. This work focuses on quantization techniques that can achieve large energy efficiency gains with minimal accuracy loss on transformers.

Prior work has also pursued optimizations specific to transformers. Algorithmic optimizations have been employed to reduce network size using techniques, such as knowledge distillation [33], inductive biases [34], and approximations [35], [36], [37], [38]. Quantization techniques with coarse-grained scaling have been used to enable INT8 for transformers [39], [40]. This work explores more aggressive quantization with INT4-VSQ while maintaining accuracy. Prior works also exploit sparsity to skip unnecessary computations in attention layers [16], [41], [42], [43], [44], and other hardware-software techniques have been explored to improve energy efficiency through early-exit prediction [45], approximate softmax [36], [38], and customized tiling schemes [46]. These techniques are complementary to our work, which also includes a hardware-efficient approximate softmax engine.

### III. ACCELERATOR DESIGN

The accelerator design focused on three key objectives: 1) high energy efficiency by exploiting low-precision computation and data reuse across different operands; 2) high area efficiency by maximizing compute throughput while amortizing memory and control overheads; and 3) programmability to support different data formats and types of computations.

#### A. Inference Accelerator Architecture

Fig. 6 shows the architecture of the accelerator, which computes a CONV, MM, or fully connected (FC) layer followed by optional post-processing operations. It consists of an array of vector MAC units, SRAM buffers to store the input matrices  $A$  and  $B$  and the output matrix  $C$ , a latch-array-based collector to accumulate partial sum values, an input register for the  $A$  operand, and a post-processing unit (PPU). The vector MAC units are configurable, performing 32-wide (INT8) or 64-wide (INT4 and INT4-VSQ) vector MAC operations each cycle, which enables direct energy efficiency comparison between the three formats and support for workloads with coarse-grained

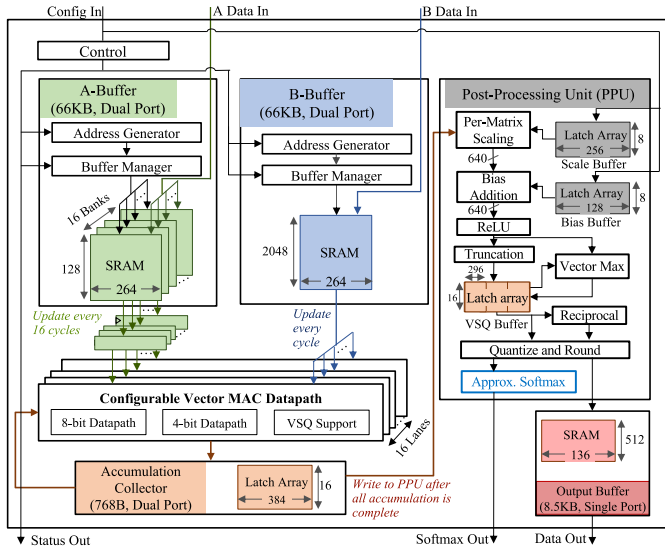


Fig. 6. Accelerator block diagram.

heterogeneous precision. The accelerator has 16 vector lanes; the 16 vector MAC units together perform a matrix-vector product and produce 16 different partial sum values each cycle that are stored in the accumulation collector.

The accelerator achieves multiple levels of reuse across all three matrices: temporal *A*-input reuse by storing vectors in the *A* register for 16 cycles, temporal output reuse through the accumulation collector, spatial reuse in the *B* input by broadcasting to each of the 16 vector MAC units, and spatial output reuse in the adder tree of vector MAC units [47]. The accumulation collector is implemented with a dual-ported latch array to enable simultaneous read/write operations every cycle at low energy cost. More details of the dataflow and mapping approach are presented in Section IV-B.

The buffers are designed to provide adequate bandwidth to the vector MAC units to achieve high throughput. The *A* buffer has 16 banks to load a different *A*-input vector for each vector lane. Each entry of the banks in the *A* and *B* buffers is 264 bits wide to store either 1) a 32-wide 8-bit vector for INT8 format; 2) a 64-wide 4-bit vector for INT4 format; or 3) a 64-wide 4-bit vector and an 8-bit scale factor for INT4-VSQ format. This matches the input width of a vector MAC unit, providing the required read bandwidth to achieve high utilization. The 16 banks of the *A* buffer supply 16 unique *A*-input vectors to each of the lanes, while the *B*-input vector read from the *B* buffer is shared across all the lanes. The input buffers *A* and *B*, which are dual-ported to support one read and one write operation per cycle per bank, employ buffer managers that enable fine-grained overlap of load and store operations in the buffers [48], avoiding the area overheads associated with double-buffering. Data are stored sequentially into these buffers, exploiting the streaming nature of DL applications. Read accesses need not be sequential; different strides, reuse patterns, and tile sizes are enabled by address generators that are programmed to generate the required sequence of addresses for read operations. The buffers enable temporal data reuse and reduce data movement cost. In particular, the partial sum values have wider widths than input operands, and their

load/store from SRAMs is expensive. Therefore, the *A* buffer and *B* buffer are sized to achieve temporal output partial sum reuse up to 128 times in the accumulation collector. This enables the accelerator to execute all of the computations in BERT-Large without requiring accesses to external memories for partial sum values.

The SRAM buffers and the latch-array collectors are shared across different data formats to improve the area efficiency of the accelerator. The width of the *A* and *B* buffers is sized so that they can optionally store per-vector scale factors alongside each vector in the INT4-VSQ format. With a bitwidth of 264, each entry in the buffers can store a 32-wide 8-bit vector, a 64-wide 4-bit vector, or a 64-wide 4-bit vector and an 8-bit per-vector scale factor. The accumulation width of each vector MAC unit is 24 bits for all the three data formats, with the accumulation collector storing 24-bit partial sum values from 16 vector lanes each cycle. All vector lanes operate in lockstep to amortize control overheads.

Once accumulation is complete, the final sum values are read from the accumulation collector to perform post-processing operations in the PPU. To ensure high utilization, the final sum reads are pipelined with the first set of store operations of the next set of partial sums. The PPU performs scaling and quantization of the partial sum values. It also supports bias addition and non-linear operations, such as ReLU and approximate softmax. The approximate softmax implementation is designed to improve hardware efficiency by using base 2 instead of base *e*, low-precision fixed-point data formats, and online normalization to reduce data movement [49]. QAT (described in IV-A) is used to mitigate the accuracy impact of the approximate softmax implementation.

The PPU is configurable to quantize the 24-bit outputs to INT8, INT4, or INT4-VSQ, depending on the desired data format for the output activations. Computing per-vector scale factors for 64-wide vectors presents certain challenges, as only 16 values are available across vector lanes and consecutive values in the accumulation collector are not ordered along the vector dimension in order to exploit weight reuse. To address this issue, the PPU includes additional latch-array buffers that store partial sum values after scaling and bias addition. Once the required values are available, the vector-max and reciprocal units compute per-vector scale factors, and the vector is quantized.

### B. Reconfigurable Datapath Design

Fig. 7 shows the microarchitecture of the single-cycle vector MAC datapath. It has dedicated datapaths that perform 32-wide INT8 and 64-wide INT4 signed dot-product computations. The output of the dot product is accumulated with signed 24-bit partial sum values for all data formats. To support INT4-VSQ, the datapath includes an unsigned 8-bit multiplier that computes the product of the two input vector scale factors. Multiplication of the 8-bit scale factors produces a 16-bit output, which is rounded to 8 bits before multiplication with the 14-bit output of the dot product. This rounding reduces the logic complexity of the datapath, improving energy efficiency with negligible impact on the accuracy of the computation.



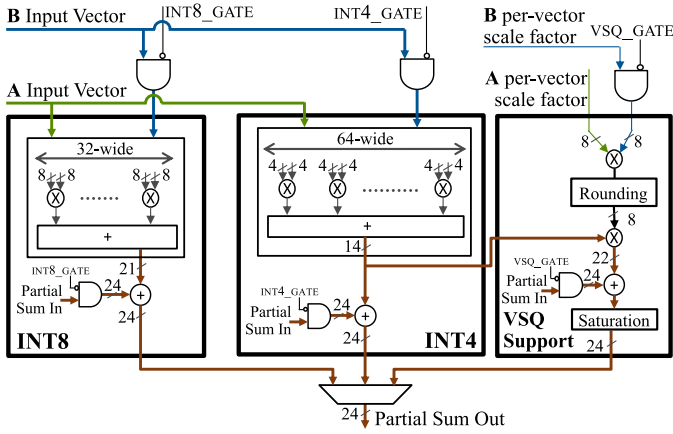


Fig. 7. Reconfigurable vector datapath.

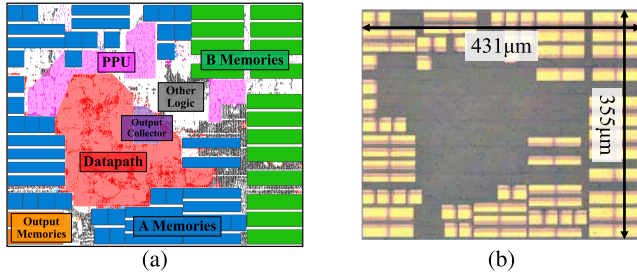


Fig. 8. (a) Annotated accelerator floorplan. (b) Die plot.

To avoid overflow errors during temporal accumulation, the datapath supports saturation of output partial sums.

The datapath can be reconfigured to support three different data formats. As shown in Fig. 7, inactive datapaths are gated by ANDing only the **B** input with zeroes, which minimizes area overhead while improving energy efficiency. For example, to configure the datapath for INT4-VSQ mode, INT4\_GATE and VSQ\_GATE are deasserted to enable the INT4 datapath and VSQ logic, while INT8\_GATE is asserted to gate the INT8 datapath.

### C. Physical Design

The physical implementation flow used to implement the accelerator favored simplicity, flexibility, and rapid turnaround time over absolute quality of results [50]. The accelerator was designed with a SystemC-based object-oriented high-level synthesis flow [51], which improved productivity by enabling design at a higher level of abstraction and speeding verification. Fig. 8(a) shows an annotated diagram of the accelerator floorplan. The entire accelerator, which is composed of 379k gates, was synthesized, placed, and routed as a single flat partition, enabling push-button SystemC-to-layout spins in <48 h. Macros were placed by the place-and-route (P&R) tool in an initial floorplanning step prior to standard cell placement. The tool was permitted to re-floorplan the macros optimally in each subsequent P&R spin, even just prior to tape out, which proved valuable in accommodating late design changes. Additional guard rings were added around each macro to ensure that arbitrary macro placements would not cause DRC or routability issues. Partition utilization was

TABLE II  
SIMULATED ACCELERATOR ENERGY BREAKDOWN

Data Format	Datapath	A-Buffer	B-Buffer	Accum. Collector	Other
INT8	2.76	0.36	0.44	0.71	1.05
INT4	1.00	0.18	0.22	0.30	0.60
INT4-VSQ	1.01	0.18	0.22	0.30	0.61

deliberately kept low to reduce P&R runtime and accommodate the macro-dominated floorplan, with a final standard cell utilization of 60.1%.

### D. Clocking

The accelerator clock is generated entirely on die via a simple multiplexor-based 64-tap ring oscillator (RO) located near the accelerator instance that is placed and routed automatically, with false-path timing constraints specified to break the combinational loop. The generated clock adapts its frequency to any changes in supply voltage, approximating the delay of the critical paths in the design. The partition is entirely synchronous except for a small amount of logic for external data transfer that is clocked by a 100-MHz reference and communicates with the accelerator via bisynchronous FIFOs. The accelerator met a clock period target of 1.1 ns at the typical 0.67-V signoff corner, with the critical paths primarily occurring from the **A** register to the accumulation collector through the INT8 datapath. The relatively relaxed frequency target improved spin times and energy efficiency while maintaining reasonable area efficiency.

### E. Energy Breakdown

Table II shows a simulated energy breakdown of the accelerator running a synthetic benchmark consisting of Gaussian random input data with 70% of the inputs set to zero (30% sparsity) in each of the supported data formats, normalized to INT4 datapath energy. Energy was measured with PrimePower at the typical 0.67-V corner using traces of glitch-aware gate-level simulations of the post-P&R netlist and normalized by the number of MAC operations performed each cycle. For all data formats, the datapath is the single largest consumer of energy, as buffer energy is reduced by amortizing accesses via the efficient dataflow described in Section IV-B and employing energy-efficient latch arrays for the shallow accumulation collectors. Control overheads are amortized by the wide vector MAC operations, and a sufficiently large tile size minimizes the relative cost of PPU computations. While some prior works claim that memory accesses dominate DL inference accelerator energy costs [52], [53], [54], [55], we demonstrate that an efficient implementation can greatly reduce non-datapath overheads.

### F. Scalable System

The described accelerator lacks a number of capabilities that would be required to accelerate neural network inference end-to-end, including a flexible high-capacity memory system, an integrated controller, and hardware support for operations such as layerwise addition, normalization, and transpose.

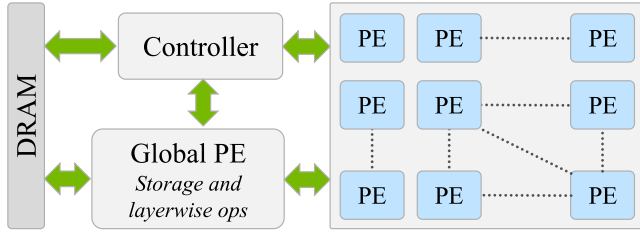


Fig. 9. Complete DL inference system composed of many instances of the proposed accelerator as processing elements (PEs).

We made the intentional decision to design an optimized accelerator core that could be scaled to a full system via replication. The core was designed to include all hardware components that would be needed for replicated cores comprising a full system. For example, the accelerator includes hardware support for streaming  $A$  and  $B$  inputs into the local buffers during execution, even though this operating mode is not supported in the silicon implementation due to the lack of a memory system. Fig. 9 shows an example of how the proposed design could be tiled into a full system.

#### IV. SOFTWARE WORKLOADS

This section describes how pretrained DNN models are fine-tuned and mapped onto the accelerator. The accelerator design was tightly coupled with the development of the software workflow.

##### A. Quantization-Aware Fine-Tuning

To achieve the best accuracy, each target model is fine-tuned from a pre-trained full-precision checkpoint targeting a specific task. Each iteration of fine-tuning consists of a hardware-aware forward pass to faithfully simulate the effects of hardware optimizations on the accuracy of the model deployed on the accelerator, followed by a backward pass based on the straight-through estimator (STE) to update the model with hardware optimizations in the loop.

1) *Forward Pass*: Tensor computations, such as CONV and MM layers, can be expressed as a set of partial sums  $S(k)$  as shown in (1), where  $I_A$  and  $I_B$  are the integer input vectors,  $V$  is the vector size,  $J$  is the number of vectors in the tensor, and  $s_A$  and  $s_B$  are the per-vector scale factors

$$S(k) = \sum_{j=0}^{J-1} \left( \sum_{i=0}^{V-1} I_A(j, i) I_B(j, i) \right) s_A(j) s_B(j). \quad (1)$$

For the INT8 and INT4 data formats,  $s_A(j) = s_B(j) = 1$ . For INT4-VSQ, the scale factors are computed using a two-level quantization scheme [21]. The first quantization level computes per-vector floating-point scale factors, and the second level quantizes the per-vector scale factors into integer per-vector components and a floating-point per-tensor component.

While a hardware-agnostic tensor-level forward pass could serve as a good baseline for abstracting the computations in an accelerator, it is insufficient for deriving the actual accuracy realized by the accelerator. As shown in Fig. 7, the per-vector

scale factor product is rounded from 16 to 8 bits prior to scaling the dot product. In addition, the partial sum saturates at 24 bits, even though its theoretical maximum exceeds this bitwidth. Moreover, fixed-point math with finite precision is used in the reciprocal operation of the PPU shown in Fig. 6 to dynamically compute the per-vector scale factors. Considering these optimizations, the actual partial sums on the accelerator  $S'(k) = S'(k, J)$  are expressed recursively as in (2), where  $\mathcal{R}$  denotes rounding and  $\mathcal{C}$  denotes saturation

$$S'(k, j_p) = \mathcal{C} \left[ S(k, j_{p-1}) + \left( \sum_{i=0}^{V-1} I_A(j_p, i) I_B(j_p, i) \right) \times \mathcal{R}[\mathcal{R}[s_A(j_p)] \mathcal{R}[s_B(j_p)]] \right]. \quad (2)$$

Accurate modeling of the accelerator requires a forward pass that represents all elements of (2). For this purpose, our hardware-aware forward pass splits the tensor-level computations into vector-level computations with attached intermediate rounding and saturation operations. Input tensors are divided into many 64-element chunks, transforming a single tensor multiplication into multiple parallel multiplications of corresponding chunks, as shown in Fig. 3. With this transformation, the result of each vector dot product is presented independently in the resulting tensors, enabling rounding and saturation to be applied on a per-vector granularity. As described in Section III-A, the accelerator uses an approximate softmax implementation and replaces the GELU operation with ReLU, so the forward pass in the QAT framework also modifies the non-linear functions to accurately capture these optimizations.

While it is necessary to divide the operations based on vector size to exactly represent the accelerator, the software model remains abstract and incurs only around  $2\times$  runtime overhead, because it does not include details such as dataflow and buffer sizes.

2) *Backward Pass*: During the backward pass, the training algorithm learns the effects of quantization, saturation, and rounding on the model using backpropagation and adapts the weights of the neural network accordingly to minimize accuracy loss. To backpropagate through these rounding and saturation operations, we apply STE [56] in the backward pass by approximating these operations as the identity function. With these approximations, it follows from (2) and (1) that  $S'(k) \approx S(k)$ , which implies that  $\nabla S'(k) \approx \nabla S(k)$ . An interpretation of this result is that since the accelerator approximates the tensor computations represented in (1), the gradients from those tensor computations can be used to train the model to target the accelerator. With this approach, QAT properly models the accelerator in the forward pass as  $S'(k)$  to correctly evaluate the actual accuracy, while reasonably approximating the gradients as  $\nabla S(k)$  in the backward pass to enable effective weight updates.  $\nabla S(k)$  can be implemented by calling well-optimized backpropagation routines of common DL packages, such as PyTorch [57], which is especially helpful in realizing performant QAT.

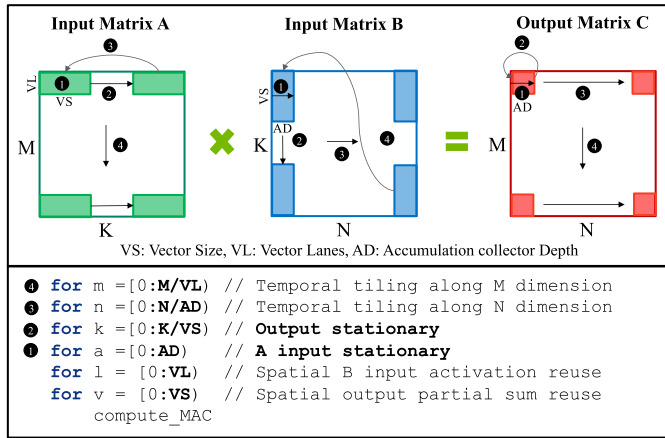


Fig. 10. Mapping MM to the accelerator.

### B. Mapping to the Accelerator

Fig. 10 shows how the different computations in deep neural networks are tiled and executed on the accelerator. In the figure,  $VS$  denotes the vector size,  $VL$  the vector lanes, and  $AD$  the accumulation collector depth. The input matrices  $A$  and  $B$  are allocated to  $A$  and  $B$  buffers, respectively.

First, matrix-vector product computation is performed by reading a sub-matrix of  $A$  of size  $VS \times VL$  and a vector of  $B$  of size  $VS$ . For INT4-VSQ format,  $VL$  per-vector scale factors of the  $A$  inputs and one per-vector scale factor of the  $B$  input are read along with the  $A$  sub-matrix and  $B$  vector. The product of the  $A$  and  $B$  per-vector scale factors is then multiplied by the matrix-vector product output within the datapaths. This produces a partial sum vector of size  $VL$ , which is stored in the accumulation collector. The  $A$  sub-matrix along with per-vector scale factors is stored in the  $A$  register and temporally reused  $AD$  times. Different  $B$  vectors are read each cycle to compute partial sum output vectors that are stored in different entries of the accumulation collector. Since the  $A$  input remains constant, it is referred to as  $A$  input stationary in Fig. 10.

Next, partial sum values in the accumulation collector are reused  $K/VS$  times to perform temporal reduction. This stage implements an output stationary dataflow by reusing partial sum outputs in the accumulation collector. This multilevel dataflow is referred to as output stationary and local  $A$  stationary, because outputs are reused in the outer loop, and  $A$  inputs are reused in the inner loop [47]. In this way, an output sub-matrix of size  $VL \times AD$  is computed in the accelerator. After the matrix-vector product, the partial sum values are quantized to the desired output format in the PPU. If the output format is INT4-VSQ,  $VS/VL$  partial sum vectors (each  $VL$  wide) are used to compute per-vector scale factors. The entire output matrix is computed via temporal tiling along the  $M$  and  $N$  dimensions. The accelerator can also efficiently execute CONV and FC layers by dividing the computation into a sequence of matrix-vector operations.

### C. End-to-End Evaluation Framework

Fig. 11 shows the end-to-end workload generation flow used to fine-tune the model and evaluate the chip. Hardware-aware

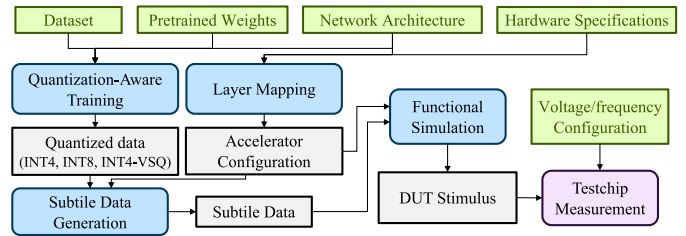


Fig. 11. Workload generation flow.

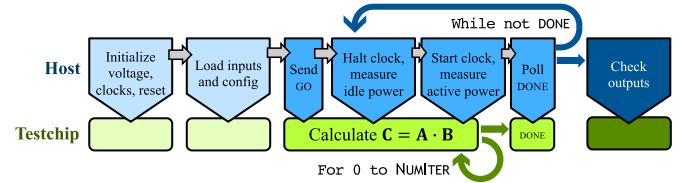


Fig. 12. Flow diagram of the chip measurement procedure.

QAT is used to recover accuracy loss from quantization and obtain quantized inputs for each layer in the desired data format. In many cases, a full network layer is too large to be executed on the accelerator, so a representative subtile of the output matrix is selected for execution. The mapping flow uses heuristics to tile the computation to exploit parallelism and data reuse in the hardware, generating the settings that configure the accelerator to perform the required computations in the vector MAC units and PPU. The quantized inputs and configuration settings are used to determine the correct data for input tiles, input scale factors, bias values, and reference outputs that represent each portion of the end-to-end DNN inference workload. This collection of data and configuration settings is used in simulation, which generates the collateral for bench measurement.

## V. MEASUREMENT RESULTS

A prototype of the accelerator was fabricated in a TSMC 5-nm process [58]. Fig. 8(b) shows a die micrograph; the testsite measures  $431 \times 355 \mu\text{m}$ .

### A. Test Setup

Fig. 12 shows the chip measurement workflow. In the initialization phase, voltages and clocks are set, and the accelerator is brought out of reset, so that its configuration registers and input memories can be programmed via the external JTAG interface for execution of a single-layer subtile. After initialization is complete, a GO command is sent to the accelerator, beginning the execution phase. Subtile execution completes in much less than a millisecond, which is an insufficient duration for accurate power measurement, and executing subsequent subtiles requires a new initialization sequence. Therefore, the accelerator state machine can be configured to execute the same subtile many millions of times in succession via a configurable NUMITER register, enabling measurement of average power. Execution termination is detected by polling a DONE register over JTAG, which is written by the accelerator once the requested number of workload iterations has been completed. In the verification phase, the outputs calculated



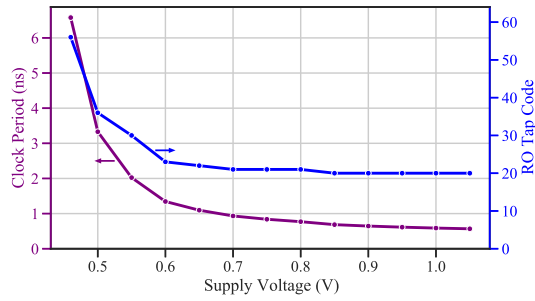


Fig. 13. Clock generator calibrated tap code and period.

by the accelerator are read out via JTAG and compared with reference outputs to confirm correct execution. Each measurement under a particular voltage and frequency condition is performed once for each subtile, with measured runtime and power extrapolated to determine the execution cost of the full tile and, hence, the full network.

The testsite supply is shared by other circuits on the die, so direct measurement of accelerator power was not possible. Instead, for a given workload, we infer the testsite dynamic power by subtracting idle power from active power and add a leakage power estimate to determine total power. To measure dynamic power, overall rail power is measured via bench measurement of board sense resistors using Agilent 34401A digital multimeters. Then, active and idle power are measured by running and then halting the accelerator clock during workload execution, with power measurements recorded during each phase (see Fig. 12). By repeating this process, five or more measurements of active and idle power are collected for each workload; the average difference between active and idle rail power is the dynamic power of the testsite for that workload. Leakage power is estimated from PrimePower reports and scaled for voltage and temperature as measured by on-die sensors. This estimate was corroborated against measured rail leakage power scaled by the area ratio of the accelerator to that of the entire voltage domain.

All measurements were performed at nominal voltage (0.67 V) unless otherwise noted. The number of operations (“ops”) used for TOPS/W and TOPS/mm<sup>2</sup> calculations is based only on the count of MAC operations, with one MAC consisting of two ops, ignoring scale factor multiplications and PPU operations.

### B. Clocking

The RO frequency is set by offline calibration of the on-die clock generator by sweeping the RO tap code and using the fastest frequency that maintains functional correctness across a representative workload subset. Fig. 13 shows the best achievable RO tap code at each voltage. Frequency is measured using integrated counters that compare against the 100-MHz reference. At voltages above 600 mV, the RO delay closely matches that of the critical paths in the design, so little adjustment in tap code is necessary to operate at peak frequency. At lower voltages, critical paths likely become dominated by wires or SRAM macros. These structures are not represented in the RO clock generator, so the number of

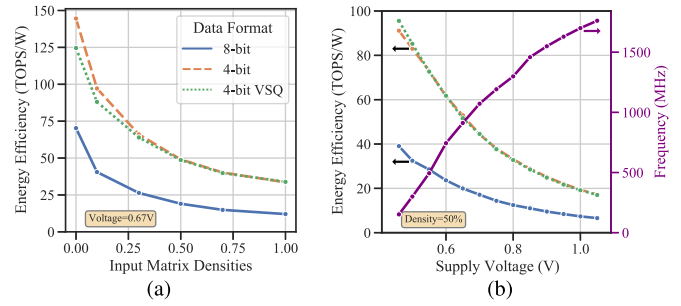


Fig. 14. Measured energy efficiency of the accelerator executing a benchmarking workload with different (a) workloads and (b) operating voltage.

multiplexors in the delay chain must be dramatically increased to track the low-voltage behavior. Each RO tap adjusts the system frequency by  $\approx 5\%$ , so a finer adjustable delay could enable further marginal performance improvement.

### C. System Benchmarking

Fig. 14 shows the measured frequency and energy efficiency of the accelerator under various operating conditions when executing a synthetic benchmark. The workload consists of an MM with **A** and **B** input values selected from a Gaussian distribution and quantized to each of the three data format modes, with values randomly zeroed out to achieve the target data density. The chip operates correctly down to a minimum voltage  $V_{\min}$  of 0.46 V and achieves a maximum frequency of 1.76 GHz at the maximum voltage  $V_{\max}$  of 1.05 V.

As Fig. 14(a) shows, the achieved energy efficiency depends not only on the choice of data format, but also on both the operating voltage and input data density. Even though the accelerator does not explicitly support sparsity through techniques, such as compressed representations or zero skipping, datapath activity is very dependent on workload sparsity, so the power consumption of the accelerator executing identical layer shapes and data format at a fixed voltage can vary by as much as  $5.9\times$ . At  $V_{\max}$  with fully dense data, the INT4-VSQ mode achieves only 11.8 TOPS/W, while at  $V_{\min}$  using 10% dense data, the accelerator achieves 156.7 TOPS/W. Using the even more unrealistic operating condition of 0% dense data (i.e., all zeroes) increases energy efficiency further to 204.5 TOPS/W. We feel that a fair comparison for benchmarking is to report energy efficiency on 50% dense data at nominal and minimum voltage. Using these criteria, our accelerator achieves 48.7 TOPS/W at 0.67 V using INT4-VSQ and 95.6 TOPS/W at 0.46 V. The use of VSQ imposes a small (0.8%) energy overhead at nominal voltage compared with baseline INT4 math.

Table III shows a comparison to prior DNN inference accelerators that support INT8 or INT4 data formats. Direct comparisons are difficult, because many prior works do not report the data density of the workloads used to measure energy efficiency. Using the reported data, our accelerator achieves a  $2.2\text{--}4.1\times$  improvement in energy efficiency and a  $3.4\text{--}15.6\times$  improvement in area efficiency for INT8 inference, and a  $5.8\times$  improvement in energy efficiency and a  $3.4\text{--}4.5\times$  improvement in area efficiency for INT4 inference.



TABLE III  
COMPARISON WITH PRIOR WORK

	This Work			Agrawal et al., ISSCC'21 [12]	Mo et al., ISSCC'21 [26]	Park et al., ISSCC'22 [23]	Zimmer et al., JSSC'21 [59]	Wang et al., ISSCC'22 [16]
Process Technology	5 nm			7 nm	28 nm	4 nm	16 nm	28 nm
Area (mm <sup>2</sup> )	0.153			19.6	1.9	4.74	6.0	6.82
Supply Voltage (V)	0.46–1.05			0.55–0.75	0.6–0.9	0.55–1.0	0.41–1.2	0.56–1.1
Frequency (MHz)	152–1760			1000–1600	100–470	332–1196	161–2001	50–510
On-Chip SRAM (kB)	141			8192	206	2048	625	336
Target Workloads	CONV, FC, Attention <sup>5</sup>			CONV, FC, RNN	CONV, FC	CNN, FC, DW-CONV	CONV, FC	Attention
Supported Data Formats	INT4	INT4-VSQ	INT8	INT2/4, FP8/16/32	INT8	INT4/8/16, FP16	INT8	INT12
Performance at $V_{\max}$ (TOPS)	3.6	3.6	1.8	102.4 (4b)	1.43	39.3 (4b), 19.7 (8b)	4.01	4.07
Energy Eff. at $V_{\min}$ (TOPS/W)	91.1 <sup>1</sup>	95.6 <sup>1</sup>	39.1 <sup>1</sup>	16.5 <sup>2</sup> (4b)	17.5 <sup>2</sup>	11.59 <sup>2</sup> (8b)	9.5 <sup>3</sup>	27.56 <sup>4</sup>
Area Eff. at $V_{\max}$ (TOPS/mm <sup>2</sup> )	23.3	23.3	11.7	5.22 (4b)	0.75	6.9 (4b), 3.45 (8b)	1.29	0.597

<sup>1</sup>50% input density. <sup>2</sup>Input density not reported. <sup>3</sup>40% input density. <sup>4</sup>90% weakly related tokens.

<sup>5</sup> Operations with low arithmetic intensity such as element-wise addition and normalization are not supported.

TABLE IV  
MEASURED PERFORMANCE ON DNN WORKLOADS

Dataset, Task	SQuAD v1.1, Reading Comprehension												ImageNet, Image Classification							
Network	BERT-Base						BERT-Large						DeiT-Small				DeiT-Base		ResNet-50	
Sequence Length	128			384			128			384			197				197		-	
Baseline FP32 Accuracy (%)	87.46			87.50			90.29			90.93			79.82				81.80		76.16	
Data Bitwidth (4V = 4b VSQ)	4	4V	8	4	4V	8	4	4V	8	4	4V	8	4	4V	8	4	4V	8	4	4V
Accuracy Loss (%)	80	0.7	0.7	81	0.5	0	88	1.1	1.1	89	0.8	0.1	29	3.6	0.7	25	1.3	0.4	0.98	0.15
MAC Utilization (%)	-	98	99	-	98	99	-	98	99	-	98	99	-	94	96	-	97	98	89	89
Throughput (inferences/s)	-	87.3	44.2	-	27.7	14.0	-	24.9	12.6	-	7.8	4.0	-	209	108	-	56.2	28.5	212	212
Energy Efficiency (inferences/s/W)	-	1734	768	-	596	281	-	513	219	-	170	80	-	3519	1451	-	1010	406	6355	4714
Energy Efficiency (TOPS/W)	-	38.7	17.2	-	42.1	19.9	-	40.4	17.3	-	42.0	19.8	-	32.0	13.2	-	35.2	14.2	51.2	38.6

### D. DNN Workload Performance

Table IV shows achieved performance, energy efficiency, and task accuracy on a variety of DNN workloads. The accelerator achieves high energy efficiency on transformers for NLP (BERT), vision transformers (DeiT), and classical CNNs (ResNet-50 [1]). For the transformer-based workloads, the reported accuracy includes the impact of hardware approximations of non-linear functions, such as softmax and GELU.

With the exception of ResNet-50, the use of INT4 results in catastrophic accuracy loss, even with QAT. INT4-VSQ achieves similar accuracy to INT8 (as well as the FP32 baseline) while achieving  $>2\times$  energy efficiency over INT8. Energy efficiencies on these workloads range from 32.0 to 42.0 TOPS/W for INT4-VSQ to 13.2 to 19.9 TOPS/W for INT8. The evaluated workloads exhibit little weight or activation sparsity, so the measured energy-efficiency results are consistent with the benchmarking measurements on dense inputs at nominal voltage.

Performance on NLP tasks is evaluated on BERT-Base and BERT-Large performing a reading comprehension task using the Stanford Question Answering Dataset (SQuAD) [60] with the sequence lengths of 128 and 384. Each task uses weights pretrained on the full English Wikipedia dataset. QAT was applied during five epochs of fine-tuning using the Adam optimizer [61] and a learning rate of  $3 \cdot 10^{-5}$ . For all transformer networks, elementwise addition and layer normalization (performed only in software) use full FP32 precision. For INT4-VSQ evaluations, FC2 layers used INT8 outputs with per-vector scale factors, which improved accuracy with negligible energy impact. All BERT workloads achieve an

TABLE V  
DATA FORMATS FOR RESNET-50 LAYERS

Data Format	Layer Name
INT8	conv1, fc
INT4	res2a_branch2a, res2a_branch1, res4{a-f}_branch2c, res5{a-c}_branch2c
INT4-VSQ	res2{a-c}_branch2{b-c}, res2{b-c}_branch2a, res3{a-d}_branch2{a-c}, res{3-4}a_branch1, res4{a-f}_branch2{a-b}, res5{a-c}_branch2{a-b}

accuracy within 1.1% of the FP32 baseline using INT4-VSQ, and BERT-Base achieves 1734 inferences/s/W with a sequence length of 128, a  $2.26\times$  energy-efficiency improvement over the 8-bit result.

Performance on CV tasks is evaluated via image classification on ImageNet [62] using both DeiT and ResNet-50. DeiT used QAT for 20 epochs, with a learning rate of  $4 \cdot 10^{-5}$  for DeiT-Small and  $10^{-5}$  for DeiT-Base. ResNet-50 was retrained using stochastic gradient descent for 30 epochs with an initial learning rate of  $10^{-4}$ , which was decayed by a factor of 10 every ten epochs, with methods in [63] used for static coarse-grained scale factor calibration and gradient backpropagation. For ResNet-50, different layers use different precision formats, as shown in Table V. The first and last layers are quantized to INT8 in both the INT4 and INT4-VSQ evaluations, as is typical in CNN quantization studies [64]. For the INT4-VSQ evaluation, layers that require large reduction length ( $K > 2048$ ) are too large to fit in the on-chip buffers while still enabling full temporal reduction of partial sums in the accumulation collector, so these layers are configured to product INT4 outputs, which is used as the data format

TABLE VI  
SPARSE BERT MEASUREMENT RESULTS

Network	BERT-Base						BERT-Large					
Sequence Length	128			384			128			384		
Weight Density (%)	67	50	33	67	50	33	67	50	33	67	50	33
Accuracy Loss (%)	-.2	.4	2.3	-.3	.4	2.6	0	.2	1.0	.2	.6	1.8
Energy Eff. Gain (%)	11	21	30	8	8	20	8	17	23	7	15	22

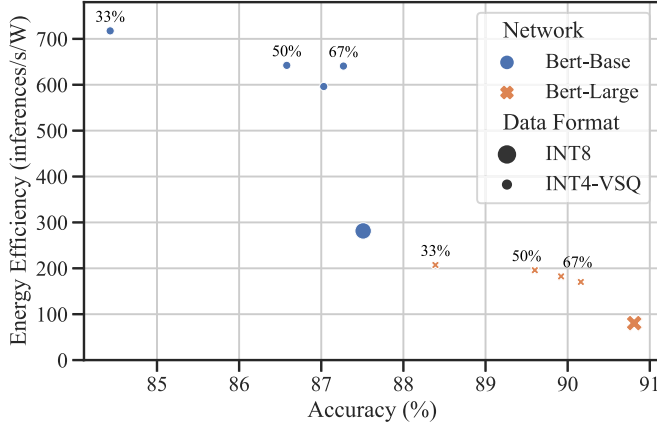


Fig. 15. Accuracy-energy trade-off for BERT with a sequence length of 384. Labeled percentages indicate weight densities after pruning.

in subsequent layers. For ResNet-50, INT4-VSQ achieves a 0.15% accuracy loss compared with the baseline, while INT4 shows a 0.98% accuracy drop.

#### E. Energy Optimizations

Software changes and hardware reconfigurability can be exploited to further improve accelerator energy efficiency. For example, while most transformer weight matrices exhibit little “natural” sparsity, weight sparsity can be increased by forcing a percentage of weights to zero during the fine-tuning process. Table VI shows the results of applying uniform per-tensor sparsity to BERT weights using magnitude-based pruning followed by retraining, similar to the approach in [65]. All accuracy and energy values are given as percent change compared with the INT4-VSQ baseline. The most aggressive sparsity (33% dense weights) results in a 20%–30% improvement in energy efficiency, though accuracy is reduced by 1.0%–2.3% due to the loss of weight information. Higher density (67% dense weights) minimizes accuracy loss, but energy efficiency gains are limited. The use of sparse training, network size, and data format can be used to generate an energy-accuracy Pareto curve (see Fig. 15).

Energy efficiency can also be improved via sparsity-aware operand ordering. Results presented thus far have mapped the  $A$  matrix into the  $A$  buffer and the  $B$  matrix into the  $B$  buffer. However, the MM transpose identity allows these operands to be reversed

$$AB = C \implies B^T A^T = C^T. \quad (3)$$

Thus,  $B^T$  can be mapped to the  $A$  buffer and  $A^T$  to the  $B$  buffer to compute a transposed version of the same result. This can save energy, because the toggle rate of the datapath

is more sensitive to the sparsity of the  $B$  input vectors, since these update every cycle, while the  $A$  input is held constant for 16 cycles in the local  $A$  stationary dataflow. Accordingly, when  $A$  is sparser than  $B$ , the transposed multiplication often costs less energy than the original. (Transpose operations are performed offline.) For INT8 BERT-Base with a sequence length of 128, layer 0 BMM2 has an  $A$  density of 98% and a  $B$  density of 64%, and energy is reduced 21% by performing the transposed multiplication with swapped operands. Additional hardware support would be required for full flexibility in operand ordering, but this preliminary result shows the promise of this simple technique when  $A$  and  $B$  have different densities.

#### VI. CONCLUSION

Energy efficiency is critical for DNN inference on transformer networks, but many prior energy-efficient DNN inference accelerators sacrifice task accuracy. This work presents an accelerator that uses VSQ to realize INT4 arithmetic for DNN inference with minimal accuracy loss on transformers. Wide vector operations and efficient dataflows that maximize reuse and minimize SRAM reads result in a prototype system that can achieve 95.6 TOPS/W in benchmarking and up to 1734 inferences/s/W running BERT on NLP workloads with <1.1% accuracy loss. This energy efficiency is 5.8× greater than the previous state of the art, setting a new benchmark for efficient DNN inference.

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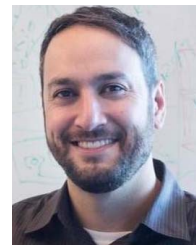


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