## FlatDD: A High-Performance Quantum Circuit Simulator using Decision Diagram and Flat Array

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## ABSTRACT

Quantum circuit simulator (QCS) is essential for designing quantum algorithms because it assists researchers in understanding how quantum operations work without access to expensive quantum computers. Traditional array-based QCSs suffer from exponential time and memory complexities. To address this problem, Decision Diagram (DD) was introduced to compress simulation data by exploring the circuit regularity. However, for irregular circuit structures, DD-based simulation incurs significant runtime and memory overhead. To overcome this challenge, we present FlatDD, a high-performance QCS that capitalizes on the strength of both DD- and array-based approaches. FlatDD parallelizes the simulation workload at multiple levels and leverages an efficient caching technique to reuse historical results. To further enhance the simulation performance for deep circuits, FlatDD introduces a gate-fusion algorithm to reduce the computational cost. Compared to state-ofthe-art QCSs on commonly used quantum circuits, FlatDD achieves 34.81× speed-up and 1.93× memory reduction.

## **CCS CONCEPTS**

• Computing methodologies  $\rightarrow$  Parallel algorithms; Simulation tools; • Computer systems organization  $\rightarrow$  Quantum computing.

### **KEYWORDS**

Quantum Circuit Simulation, Decision Diagram, Exponentially Weighted Moving Average

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## **1 INTRODUCTION**

Quantum computing (QC) has the potential to efficiently handle certain types of problems that are classically intractable, such as quantum chemistry [9], finance [35], and cryptography [20]. Powered by two fundamental quantum phenomena, *superposition* and *entanglement*, many available quantum computers have shown significant speed-up over classical computers. For example, a recent photonic quantum computer Jiuzhang can solve the Gaussian boson sampling problem within five minutes, whereas a supercomputer needs billions of years [95]. To enable widespread use of this quantum advantage, researchers have been actively building software stacks to support quantum computer designs [4, 11, 41, 81, 82, 90].

Among various QC applications, developing an efficient quantum circuit simulator (QCS) on a classical computer is a crucial task because it helps researchers understand how quantum operations work and verify the functionality of a quantum algorithm. However, this task is extremely challenging because it demands large space and time complexity to compute state amplitudes of qubits. For instance, state-of-the-art QCSs [1, 4, 5, 19, 63, 68, 89, 91, 94] use arrays to represent quantum gate matrices and state vectors. An n-qubit circuit may result in a worst case of multiplying a  $2^n \times 2^n$  quantum gate matrix by a  $2^n \times 1$  state vector. To tackle this challenge, [86, 99] have proposed decision diagram (DD) to compress simulation data in a compact graph-based data structure by exploring regularity in the circuit, such as state amplitude distribution and gate matrices structures. As a result, DD can significantly reduce the space complexity and largely improve the simulation time. It is widely used by many quantum software projects [11, 21, 22, 36, 37, 97].

Despite superior performance on regular circuit structures (e.g., quantum arithmetic), DD-based simulators cannot efficiently handle circuits that exhibit irregular distributions of state amplitudes, such as deep neural network (DNN) [10], variational quantum eigensolver (VQE), and Google's quantum supremacy circuits [7]. When simulating these irregular circuits, DD has few advantages but suffers from exponential runtime and memory overhead due to its graph structure–which would otherwise be more efficient using 1D arrays. Figure 1 illustrates this problem by showing the runtime and memory results between a DD-based simulator [99] and an array-based simulator [19] on two regular (Adder, GHZ State [88]) and two irregular (DNN, VQE) quantum circuits.

To solve this problem, we have identified an important property: In DD-based simulation, the quantum gate matrix has a regular



Figure 1: Normalized runtime and memory results between a DD-based simulator [99] and an array-based simulator [19] on four structurally different quantum circuits.

structure as it can be recursively decomposed to unitary operators through Kronecker product. On the other hand, quantum state vectors exhibit a different behavior because of superposition and amplification over the state space [21]. Typically, the vector starts with a highly regular distribution and gradually becomes irregular as the simulation progresses. With this property, we present a highperformance QCS called FlatDD that capitalizes on the strength of both DD- and array-based approaches. We summarize our technical contributions below:

- We introduce a hybrid data structure *DMAV*, which uses <u>DD</u>based gate <u>matrix</u> and <u>array-based</u> state <u>vector</u> for matrix-vector multiplication. DD-based gate matrix enhances indexing efficiency, while array-based state vector avoids exponential overhead from irregularity.
- We introduce a parallel DMAV algorithm with an efficient caching technique to reuse simulation data. Our DMAV algorithm overcomes the limitation of DD-based simulation which is inherently sequential, thus largely enhancing its runtime scalability on a multicore architecture.
- We introduce a moving average-based algorithm to effectively decide when to convert the simulation from DD to DMAV. Since such a conversion can be time-consuming for large circuits, we introduce a parallel algorithm to maximize the conversion efficiency.
- We introduce a DMAV-aware gate-fusion algorithm to enhance FlatDD's efficiency in handling large quantum circuits with deep simulation length.

We evaluated FlatDD on a set of widely-used quantum circuits from [7, 69, 88]. FlatDD can outperform two highly optimized stateof-the-art DD-based and array-based simulators, DDSIM [99] and Quantum++ [19], with significant runtime improvement. For example, FlatDD achieves an average of 34.81× and 17.31× speed-up over DDSIM [99] and Quantum++ [19]. The source code is available at https://github.com/IDEA-CUHK/FlatDD.

#### 2 QUANTUM CIRCUIT SIMULATION

Given a quantum circuit, the goal of quantum circuit simulation is to derive the final state value after applying all quantum gate operations to an initial state. Mainstream simulation methods can be categorized to array-based and DD-based, explained below:



Figure 2: Examples of array-based data represented in decision diagrams. Edges without labels have a weight of one by default.

#### 2.1 Array-based Simulation

In array-based simulation [1, 4, 5, 19, 63, 68, 89, 91, 94], gate matrices and state vectors are stored in 2D and 1D arrays. Multiplying a 2D array with a 1D array expresses the action of exerting a quantum gate on a state vector. For instance, when we apply a single-qubit Hadamard gate *H* to an input state  $|\psi\rangle = |0\rangle$ , it yields the resulting state  $|\psi'\rangle$  (Equation 1).

$$|\psi'\rangle = H \cdot |\psi\rangle = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1\\ 1 & -1 \end{pmatrix} \cdot \begin{pmatrix} 1\\ 0 \end{pmatrix} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1\\ 1 \end{pmatrix}$$
(1)

This can also extend to larger circuits with n > 2 qubits. However, we do not have to construct an entire  $2^n \times 2^n$  gate matrix [89, 91]. Instead, array-based simulators can manipulate the amplitudes of state vectors locally. For instance, if we apply a single-qubit gate  $U = (u_{ij})$  to the *k*-th qubit of state vector  $|\psi\rangle = (a_i)^T$ , the operation to derive resulting state  $|\psi'\rangle = (a'_i)^T$  can be expressed in Equation 2.

$$\begin{pmatrix} a'_{*\cdots*0_{k}*\cdots*} \\ a'_{*\cdots*1_{k}*\cdots*} \end{pmatrix} = \begin{pmatrix} u_{11} & u_{12} \\ u_{21} & u_{22} \end{pmatrix} \cdot \begin{pmatrix} a_{*\cdots*0_{k}*\cdots*} \\ a_{*\cdots*1_{k}*\cdots*} \end{pmatrix}$$
(2)

On the other hand, if we apply a two-qubit controlled gate  $V = (v_{ij})$  to the state vector, where *c* represents the control qubit and *t* is the target qubit, the in-place manipulation of the state vector is expressed in Equation 3.

$$\begin{pmatrix} a'_{*1_c*\cdots*0_t*\cdots*}\\ a'_{*1_c*\cdots*1_t*\cdots*} \end{pmatrix} = \begin{pmatrix} v_{11} & v_{12}\\ v_{21} & v_{22} \end{pmatrix} \cdot \begin{pmatrix} a_{*1_c*\cdots*0_t*\cdots*}\\ a_{*1_c*\cdots*1_t*\cdots*} \end{pmatrix}$$
(3)

#### 2.2 Decision-diagram-based Simulation

Compared with array-based methods, decision diagram (DD)-based simulators [86, 98, 99] are particularly good at handling the regularity in gate matrices and state vectors. For example, if we equally partition the matrix in Figure 2a into four sub-matrices, we observe that the upper-left, upper-right, and lower-left sub-matrices are identical. The lower-right sub-matrix is exactly the opposite of the other three sub-matrices. As a result, these four sub-matrices can be efficiently stored as one single  $2 \times 2$  sub-matrix with different weights. This idea extends to all gate matrices and state vectors, which can be recursively partitioned until they become scalar values. This process forms a graph-based data structure, DD, where nodes represent sub-matrices or sub-vectors at various partitioning levels connected by weighted edges. Figure 2 shows DD examples for gate matrix *M* and state vector *V*.

In Figure 2a, the root node  $m_1$  represents the entire matrix M, and the four outgoing edges represent four equally partitioned submatrices in M. The partition runs on the most significant qubit,  $q_1$ , at the topmost level of the DD. Recursive partitions progress level by level, from the most to the least significant qubit. Each qubit matches a unique level in DD. The weight for  $m_1$ 's incoming edge is  $\frac{1}{\sqrt{2}}$ , and the weights for its outgoing edges are 1, 1, 1 and -1, corresponding to the upper-left, upper-right, lower-left, and lower-right sub-matrices. The weights are uniquely decided by normalization [86, 99]. After dividing  $m_1$ 's incoming and outgoing weights, the four sub-matrices are identical, which can be expressed with one single node  $m_2$ .  $m_2$  represents a 2 × 2 identity matrix, which can be further partitioned on qubit  $q_0$ . In this partition, the upper-left and lower-right elements point to terminal constant node one, and the upper-right and lower-left elements are zero. The matrix value of an index pair equals the product of edge weights along the corresponding path in DD. For instance, for M[0][2] (i.e.,  $M[|00\rangle][|10\rangle])$ , we have  $q_1 = q_0 = 0$  and  $q_1 = 1$ ,  $q_0 = 0$  for row and column indices, respectively. Multiplying the edge weights along the path (thick red edges in Figure 2a) yields  $M[0][2] = \frac{1}{\sqrt{2}} \times 1 \times 1 =$  $\frac{1}{\sqrt{2}}$ .

In Figure 2b, the root node  $v_1$  represents the entire state vector. We equally partition the vector into two sub-vectors on the most significant qubit  $q_2$ . The two sub-vectors are neither zero vectors nor a scalar multiple of each other. Therefore, they are represented in two unique nodes  $v_2$  and  $v_3$ . The incoming weights for  $v_2$  and  $v_3$  are  $\frac{1}{\sqrt{2}}$ , also uniquely determined by normalization. Thus, the weight products along the paths to  $v_2$  and  $v_3$  are both  $\frac{1}{\sqrt{2}}$ . After dividing the weight products, the two sub-vectors represented by  $v_2$ and  $v_3$  can be further partitioned into  $(\frac{1}{\sqrt{2}} \ 0)^T$ ,  $(0 \ \frac{1}{\sqrt{2}})^T$ ,  $(\frac{1}{\sqrt{2}} \ 0)^T$ and  $\left(0 - \frac{1}{\sqrt{2}}\right)^T$  on  $q_1$  level, where the first and the third are identical, represented in node  $v_4$ , and the second and the fourth are opposite, represented in node  $v_5$  with opposite incoming weights. After further dividing weight products,  $v_4$  and  $v_5$  represent  $\begin{pmatrix} 1 & 0 \end{pmatrix}^T$  and  $\begin{pmatrix} 0 & 1 \end{pmatrix}^T$  on the  $q_0$  level, respectively. Similarly, the amplitude of an index is equal to the product of edge weights along the corresponding path in DD. For example, to determine V[3] (i.e.,  $V[|011\rangle]$ ), we have  $q_2 = 0, q_1 = 1$  and  $q_0 = 1$ . Multiplying the edge weights along the path (thick red edges in Figure 2b) gives  $V[3] = 1 \times \frac{1}{\sqrt{2}} \times \frac{1}{\sqrt{2}} \times 1 = \frac{1}{2}$ .

DD-based matrix-vector multiplication is done in a depth-firstsearch (DFS) fashion, where each matrix node always finds its corresponding vector node counterpart on the same level. Identical matrix-vector multiplications are avoided using hash tables [86, 98, 99].

### **3 ALGORITHM**

In this section, we discuss the technical details of our FlatDD algorithm. Figure 3 gives an overview of FlatDD. As aforementioned, the quantum state vector typically begins with a highly regular distribution and gradually turns irregular as the simulation progresses. For example, in Figure 3, DD-based simulation has a fast runtime until about 24 gates at which the cost to maintain an irregular DD-based state vector outweighs its advantage. Considering this property, FlatDD begins with DD-based simulation using DDSIM [99] and converts to parallel simulation with DD-based gate matrix and array-based state vector multiplication (DMAV). We



Figure 3: Overview of the FlatDD algorithm. The top box displays the runtime for each quantum gate, while the bottom box shows the quantum circuit and simulation data structures used to compute state vectors.

record a window of the DD size using moving average and examine if the state vector's regularity has experienced a significant change. When such a change happens, we use a parallel algorithm to convert the state vector from DD to array (Section 3.1), thereby converting the simulation from DD-based to DMAV. In Section 3.2, we introduce an efficient parallel DMAV algorithm with a caching technique to maximize performance. To further enhance the performance of FlatDD on large, deep circuits with thousands of gates, we introduce a DMAV-aware gate-fusion algorithm (Section 3.3). Throughout this paper, we use *t* to represent the number of threads and *n* to represent the number of qubits.

## 3.1 Conversion from DD-based Simulation to DMAV

When the regularity of a DD-based state vector decreases to a certain level, DD-based simulation becomes less advantageous due to the cost of maintaining a highly irregular DD structure. To address this problem, a potential solution is to convert the DD-based state vector to an array representation when the DD size exceeds a certain threshold. The challenge is thus to decide when and how to perform this conversion. To this end, we decide the conversion timing via moving average (Section 3.1.1) and introduce a parallel algorithm to efficiently convert DD to array (Section 3.1.2).

3.1.1 Conversion timing. To decide the best timing for conversion from DD to array, we monitor the DD size and perform conversion when the regularity of DD experiences a drastic increase. Specifically, we apply a widely used learning method, *exponentially weighted moving average* [59] (EWMA). EWMA introduces little computational overhead and is suitable for different quantum circuits and simulation platforms. While simulating, gate *i* is assigned an EWMA value  $v_i$  based on the previous EWMA value  $v_{i-1}$  and the DD size  $s_i$ , of the state vector at gate *i*.  $v_i$  is computed through Equation 4.

$$v_i = \beta \cdot v_{i-1} + (1 - \beta) \cdot s_i \tag{4}$$

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(a) Load balancing: Avoiding idle (b) Scalar multiplication: Facilithreads at zeros. tating SIMD parallelism.

Figure 4: Parallel DD-to-array conversion algorithm with two optimizations. The four colors represent four threads.

In Equation 4,  $\beta$  is a parameter and  $v_0$  is initialized to 0.  $\beta$  captures the learning feature of EWMA based on a weighted window of historical data. To decide whether to convert at gate *i*, we compare  $\epsilon \cdot v_i$  with  $s_i$ , where  $\epsilon$  is a threshold parameter. (1) If  $\epsilon \cdot v_i \ge s_i$ , then  $s_i$  is not large enough to benefit from DMAV. (2) On the other hand, if  $\epsilon \cdot v_i < s_i$ , we convert from DD-based simulation to DMAV (Section 3.2) by converting state vector from DD to 1D flat array.

3.1.2 Parallel Conversion from DD to Array. Although DD-based QCS DDSIM [99] already incorporates a conversion algorithm, it is inherently sequential. As a result, it can consume a large portion of the total simulation runtime. To overcome this challenge, we introduce a parallel algorithm with two optimizations, load balancing and scalar multiplication, to improve the conversion efficiency.

Figure 4 illustrates how our parallel conversion algorithm converts state vector V from DD to array using four threads  $(t_1, t_2, t_3, t_3, t_4)$ and  $t_4$ ). In our algorithm, each DD node serves as a junction that divides threads across its two outgoing edges, progressing from top to bottom until further division of threads is not possible. Then, each thread traverses the unvisited nodes using depth-first-search (DFS), and computes the state amplitudes through the products of weights along the traversal paths. In practice, however, edges may be zero. For example, as shown in Figure 4a, the right outgoing edges of  $v_1$  and  $v_3$ , and the left outgoing edge of  $v_4$  are zero. We address this problem with load-balancing optimization: If a DD node's outgoing edge is zero, threads are not divided; instead, they all proceed along the non-zero edge. With this optimization, in Figure 4a, all four threads follow the left outgoing edge at node  $v_1$ to node  $v_2$ . Here, the threads divide: two take left to  $v_3$  and two take right to  $v_4$ . At nodes  $v_3$  and  $v_4$ , the threads continue to nodes  $v_5$  and  $v_6$ , respectively, and divide at  $v_5$  and  $v_6$ . This optimization prevents any threads from remaining idle.

Additionally, we optimize the conversion by identifying opportunities for scalar multiplication, as it benefits from single instruction multiple data (SIMD) parallelism. For example, in Figure 4b, we can derive the second, third, and fourth quarters of the array from its first quarter with scalar multiplication. This is because  $v_1$  and  $v_2$ each have identical child nodes, indicating that the four quarters of the array are scalar multiples of one another [86]. Consequently, we break the conversion in Figure 4b into two steps: ① Convert the first quarter. With multi-threading, all four threads make two consecutive left turns at  $v_1$  and  $v_2$ , for they each have identical child nodes, and divide at nodes  $v_3$ ,  $v_4$ , and  $v_5$ , to convert the first quarter in parallel. ② SIMD-enabled scalar multiplication fills the second, third, and fourth quarters with three threads, using the first quarter data. For instance, the fourth quarter is calculated by multiplying the first quarter by scalar bd/ac. This optimization enhances performance by leveraging both SIMD and multi-threading parallelism.

#### 3.2 Simulation with DMAV

DMAV is different from existing matrix-vector multiplications [8, 85] because a DD-based gate matrix has a specific regularity property due to the tensor product. This regularity allows reusing computed results through caching. However, caching requires additional memory to store historical data, which can introduce overhead that, for certain gates, can outweigh its benefits. To effectively determine which gates benefit from caching, we introduce a computational cost model for DMAV (Section 3.2.3). Using this model, we apply DMAV without caching (Section 3.2.1) when its computational cost is lower. Otherwise, we apply DMAV with caching (Section 3.2.2).

3.2.1 DMAV without Caching. We describe DMAV without caching, using Algorithm 1 and an example in Figure 5. In Figure 5, we multiply a three-qubit (i.e., n = 3) gate matrix M by a state vector V using two threads (in blue and red, t = 2), to derive state vector W. Algorithm 1 consists of three functions: DMAV, Assign, and Run. The top-level function DMAV takes input variables M (the topmost edge to DD node  $m_1$  in Figure 5), as well as V and W (the input and output state vectors in flat arrays). DMAV has two steps: assigning multiplication tasks to be executed on t threads via Assign (line 2), and running them in parallel via Run (lines 3-5).

In DMAV without caching, each thread evaluates M in row space by computing h rows from M and the entire vector V, resulting in an *h*-sized sub-vector in *W*, where  $h = 2^n/t$ . Assign decomposes this process for each thread into smaller multiplication tasks. Assign divides matrix *M* into  $h \times h$  sub-matrices, and vectors *V* and W into h-sized sub-vectors. Each sub-matrix is paired with the corresponding sub-vectors, forming a multiplication task. A DD sub-matrix can be located using its incoming edge, while a subvector can be located using its start index in V or W. Thus, to record the multiplication tasks for t threads, we use 2D vectors  $v_M, v_V, v_f$ (each of length *t*) to track the sub-matrices' DD edges, sub-vectors' start indices, and the weight products along the DD traversal paths. Assign is a recursive function, where the input arguments are dynamically decided during the recursive call. At line 8,  $M_r$  is the input sub-matrix's DD edge, and  $f_r$  is the weight product along the DD traversal path. If each DD edge  $M_{r_i}$  has weight  $M_{r_i}$ .w along DD traversal path *P*,  $f_r$  is given by  $f_r = \prod_{i \in P} M_{r_i}$ .w. Argument *u* is the thread index for task assignment.  $I_V$  is the start index for a sub-vector in V. l denotes the DD level of the current node, pointed to by edge  $M_r$  (i.e.,  $M_r.n$ ). Assign is called from DMAV (line 2) using DD edge *M*. The call initializes weight product  $f_r$ , thread index *u*, and start index  $I_V$  to 1, 0, and 0 respectively, on the topmost level l = n - 1.

In Assign, if  $M_r$  is zero, it returns (line 9).  $n - \log_2 t - 1$  represents the *border level*: Assign ends here, and Run starts from here. In Figure 5, the border level is  $q_1$  (i.e.,  $n - \log_2 t - 1 = 1$ ). Assign spans levels  $q_2$  and  $q_1$  (i.e., l = 2, 1), while Run spans levels  $q_1, q_0$ , and

the final level where the bottom-most terminal node one resides (i.e., l = 1, 0, -1). Upon reaching the border level, we push the submatrix's DD edge  $M_r$ , the sub-vector's start index  $I_V$ , and the weight product  $f_r$  of thread u to the corresponding vectors (lines 10-11). If the border level is still not reached, we traverse the four outgoing edges of  $M_r.n$  in row-major order (i.e.,  $M_r.n.e[i][j], \forall i, j \in \{0, 1\}^2$ ) and call Assign with the updated weight product  $M_r.w \cdot f_r$  (where  $M_r$ . w is the weight of  $M_r$ ), thread index  $u+i\cdot t/2^{n-l}$  and sub-vector's start index,  $I_V + 2^l j$ , going one level lower to l - 1 (lines 12-13). The purposes of  $u + i \cdot t/2^{n-l}$  and  $I_V + 2^l j$  are to split *t* threads and V into halves, quarters, and so on, at each successive level, until all threads are allocated a multiplication task. In Figure 5, on level  $q_2$ , we call Assign with u = 0, 1 and  $I_V = 0, 4$ , essentially halving t threads and V. After Assign, the blue thread in Figure 5 is assigned  $a \cdot m_2 \cdot V[0:4]$  and  $b \cdot m_2 \cdot V[4:8]$ , while the red thread is assigned  $c \cdot m_2 \cdot V[0:4]$  and  $d \cdot m_2 \cdot V[4:8]$ .

Algorithm 1 DMAV without caching

**Input:**  $h = \frac{2^n}{t}$ , the size for a sub-matrix or a sub-vector;  $v_M = v_V = v_f = \{\{\}, \dots, \{\}\}$ , vectors keeping track of the sub-matrices' DD edges, sub-vectors' start indices, and weight products as-

signed to each thread, respectively. 1: **function** DMAV(M, V, W)

Assign(M, 1, 0, 0, n - 1)2: **parallel for**  $i \in [0, t)$ 3: for  $j \in [0, size(v_M[i]))$ 4:  $\operatorname{Run}(v_{M}[i][j], V, W, n - \log_{2} t - 1, v_{V}[i][j], ih, v_{f}[i][j])$ 5: end function 6: 7: 8: **function** Assign $(M_r, f_r, u, I_V, l)$ 

if M<sub>r</sub> is zero edge then return 9:

```
if l == n - \log_2 t - 1 then
10:
```

```
v_M[u] \cup \{M_r\}; v_V[u] \cup \{I_V\}; v_f[u] \cup \{f_r\}; return
11:
```

```
12:
```

```
else for i, j \in \{0, 1\}^2
Assign(M_r.n.e[i][j], M_r.w \cdot f_r, u + \frac{i \cdot t}{2^{n-l}}, I_V + 2^l j, l - 1)
13:
14: end function
```

```
15:
```

16: **function** Run( $M_r$ , V, W, l,  $I_V$ ,  $I_W$ ,  $f_r$ ) if  $M_r$  is zero edge then return 17:

```
if M_r.n is terminal node then
18:
```

```
W[I_W] \leftarrow W[I_W] + f_r \cdot M_r.w \cdot V[I_V]; return
19:
```

```
for i, j \in \{0, 1\}^2
20:
```

 $Run(M_r.n.e[i][j], V, W, l - 1, I_V + 2^l j, I_W + 2^l i, f_r M_r.w)$ 21: 22: end function

Run is also a recursive function. We launch t threads (line 3), and iterate through the edges in  $v_M$  determined by Assign (line 4). At line 5, we calculate each multiplication task  $v_f[i][j] \cdot v_M[i][j]$ .  $V[v_V[i][j] : v_V[i][j] + h]$  via Run. The input parameters (line 16) for Run consist of the following:  $M_r$ , sub-matrix's DD edge; V and W, the input and output state vectors; l, the level on which node  $M_r.n$  is located;  $I_V$  and  $I_W$ , start indices for accessing V and W; and  $f_r$ , weight product. In Run, if  $M_r$  is zero, then we return directly (line 17). If M<sub>r</sub>.n is a terminal node (lines 18-19), we multiply



Figure 5: DMAV without caching on two threads, illustrated in blue and red colors. The input matrix and vector are M and V, and the resulting vector is W (i.e.,  $M \cdot V \rightarrow W$ ).

 $M[I_W][I_V]$  by  $V[I_V]$ , and add the result to  $W[I_W]$ . If  $M_r.n$  is not yet terminal, we traverse the four outgoing edges of  $M_r.n$  and call Run accordingly. We update the weight product with  $f_r \cdot M_r.w$ ,  $I_V$ with  $I_V + 2^l j$ , and  $I_W$  with  $I_W + 2^l i$  (lines 20-21). The purposes of  $I_V + 2^l j$ ,  $I_W + 2^l i$  are to divide the *h*-sized sub-vectors in V and W into halves, quarters, and so on, at each successive level, until the sub-vectors cannot be divided further, at which point  $M_r$  also reaches a terminal node.

In Figure 5, the blue thread first computes  $a \cdot m_2 \cdot V[0:4]$ (allocated by Assign). We make two recursive Run calls, using the upper-left outgoing edges of  $m_2$  and  $m_3$  for the  $M_r$  inputs, and setting both  $I_V$  and  $I_W$  to 0, leading to  $W[0] \leftarrow W[0] + M[0][0]$ . V[0] at step (1). Then, at step (2), we exit the outermost Run and call another Run ( $M_r = m_3$ 's lower-right edge, and  $I_V = I_W = 1$ ), leading to  $W[1] \leftarrow W[1] + M[1][1] \cdot V[1]$ . This continues until  $a \cdot m_2 \cdot V[0:4]$  is finished and we make a switch for task  $b \cdot m_2 \cdot V[4:$ 8] at step (5). A similar process applies to the red thread.

DMAV outperforms array-based QCSs (e.g., Quantum++ [19]) due to its efficient indexing. Unlike Quantum++, which requires O(n)operations per state, DMAV uses a recursive Run function within a DD structure, enhancing data locality and reducing indexing to a constant average number of operations. This results in an  $n \times$ increase in indexing speed for DMAV compared to Quantum++.

3.2.2 DMAV with Caching. Figure 6 illustrates an example of computation reduction with caching in DMAV. Executing the DMAV in Figure 6 involves two multiplications: (1)  $am_2 \cdot V[0:2]$  and (2)  $bm_2 \cdot V[0:2]$ . Without caching, both require four multiplyaccumulate (MAC) [2] operations. However, (1) and (2) share identical left and right operands (sub-matrix DD node and sub-vector) and differ only in their scalar coefficients (a or b). Caching the result of (1) allows its reuse for (2) by scaling it by b/a, yielding only two MAC operations, compared to four without caching.

DMAV with caching is notably different from DMAV without caching. Specifically, to facilitate data reuse, each thread evaluates the gate matrix in column space instead of row space. We describe DMAV with caching using both Algorithm 2 and Figure 7. In Figure 7, we multiply gate matrix M by state vector V using four threads, each with its own cache, to obtain the resulting state vector W. In Algorithm 2, we present two functions: top-level DMAVCache,



# Figure 6: Saving the number of MAC operations by reusing historical results in DMAV.

which conducts DMAV with caching, and AssignCache, which assigns multiplication tasks to *t* threads. For previously computed multiplication tasks, we can fetch the results from the cache, while other multiplication tasks are computed using Run from Algorithm 1. We only apply caching to the multiplication tasks directly assigned by AssignCache to avoid frequent memory allocation and deallocation on different threads.

In DMAV with caching, each thread computes h columns in M and an h-sized sub-vector in V ( $h = 2^n/t$ ), generating a partial output equivalent in size to V. These partial outputs are then summed to obtain the output state vector W. Given the sparse nature of quantum gate matrices [34], the partial outputs from different threads often have non-overlapping segments. Thus, to save memory and time, multiple partial outputs can share a single memory *buffer*, and we sum the buffers to obtain the output state vector.

AssignCache, which is adapted from Assign in Algorithm 1, divides matrix M into  $h \times h$  sub-matrices, and vector V as well as the partial output vectors into h-sized sub-vectors. Each submatrix is paired with the corresponding sub-vectors, forming a multiplication task. To record the multiplication tasks for t threads, we use vectors  $v_M$ ,  $v_P$ , and  $v_f$  to keep track of the sub-matrices' DD edges, start indices of sub-vectors in a partial output, and weight products assigned to each thread. Vector  $v_B$  associates each thread with a specific buffer index, which can be used to access buffers in *B*, the vector of buffers. The input parameters for AssignCache are:  $M_r$ , input matrix's DD edge;  $f_r$ , weight product along DD traversal path; u, thread index;  $I_P$ , start index of a sub-vector in a partial output; and *l*, DD level of the current node. AssignCache returns if  $M_r$  is zero (line 17). Upon reaching the border level at  $n - \log_2 t - 1$ , we push the sub-matrix's edge  $M_r$ , the sub-vector's start index  $I_P$ , and the weight product  $f_r$  of thread u to the corresponding vectors (lines 18-19). At lines 20-21, we traverse the four outgoing edges of  $M_r.n$  in column-major order and call AssignCache with the updated weight product, thread index, sub-vector's start index, and DD level, similar to Assign (line 13 in Algorithm 1). After all threads have been assigned multiplication tasks, we assign each thread to its buffer. Specifically, for each thread *i*, we check if there is another thread *j* with a non-overlapping partial output. If such a thread *j* exists, threads *i* and *j* will share a buffer (line 24); otherwise, thread *i* will receive its own buffer (line 25).

In DMAVCache, each thread executes its multiplication tasks from AssignCache (line 2). We focus on threads  $t_1$  and  $t_2$  in Figure 7 (in black dashed boxes).  $t_1$  is assigned  $a \cdot m_4 \cdot V[0:h]$  and  $b \cdot m_4 \cdot V[0:h]$ , while  $t_2$  is assigned  $c \cdot m_5 \cdot V[h:2h]$  and  $d \cdot m_5 \cdot V[h:2h]$ , where sub-vectors V[0:h] and V[h:2h] are the first two quarters in V. As these four tasks have non-overlapping outputs, they can Algorithm 2 DMAV with caching

**Input:** 
$$h = \frac{2^n}{t}$$
, the size for a sub-matrix or a sub-vector;  $v_M = v_P = v_f = \{\{\}, \dots, \{\}\}\}$ , vectors keeping track of the sub-matrices?

DD edges, start indices of sub-vectors in a partial output, and weight products, respectively;  $v_B$ , a vector of buffer indices assigned to each thread; B, a vector of buffers for partial outputs; *cache*, a vector where each element is a cache specific to a thread.

1: **function** DMAVCache(M, V, W)

- 2: AssignCache(M, 1, 0, 0, n 1)
- 3: **parallel for**  $i \in [0, t)$

```
4: for j \in [0, size(v_M[i]))
```

5:  $r \leftarrow cache[i].lookup(v_M[i][j].n)$ 

```
6: if r \neq \{\} then
```

7:  $B[v_B[i]][v_P[i][j] : v_P[i][j]+h] \leftarrow \text{SIMDMul}(v_f[i][j]/r[0], B[v_B[i]][r[1] : r[1]+h])$ 

```
8: else
```

15

9:  $\operatorname{Run}(v_{M}[i][j], V, B[v_{B}[i]], n - \log_{2} t - 1, h \cdot i, v_{P}[i][j], v_{f}[i][j]) /* \operatorname{From Algorithm 1 */}$ 

```
10: cache[i].insert(v_M[i][j].n, \{v_f[i][j], v_P[i][j]\})
```

```
11: parallel for i \in [0, t)
```

```
12: for j \in [0, size(B))
```

```
13: W[i \cdot h : (i+1) \cdot h] \leftarrow \text{SIMDAdd}(W[i \cdot h : (i+1) \cdot h], B[j][i \cdot h : (i+1) \cdot h])
```

```
14: end function
```

```
16: function AssignCache(M_r, f_r, u, I_P, l)
```

```
17: if M_r is zero edge then return
```

```
18: if l == n - \log_2 t - 1 then /* Border level */
```

19:  $v_M[u] \cup \{M_r\}; v_P[u] \cup \{I_P\}; v_f[u] \cup \{f_r\};$  return

```
20: else for j, i \in \{0, 1\}^2
```

21: AssignCache( $M_r.n.e[i][j], f_rM_r.w, u + \frac{j \cdot t}{2^{n-l}}, I_P + 2^l i, l-1$ )

```
22: if l == n - 1 then for i \in [0, size(v_P))
```

```
if not Overlap(v_P[i], v_P[j]), \exists j \in [0, i) then
```

```
24: v_B[i] \leftarrow v_B[j]
```

```
25: else v_B[i] \leftarrow size(B); B \cup \{\{0, \cdots, 0\}\}
```

26: end function

share the same output buffer B[0]. At step ①,  $t_1$  and  $t_2$  execute  $am_4 \cdot V[0:h]$  and  $cm_5 \cdot V[h:2h]$  simultaneously. First, they search in caches (line 5). Finding no historical results, they directly call the Run function from Algorithm 1. Upon obtaining the result,  $t_1$  caches  $m_4$  with a pair consisting of  $m_4$ 's weight product and the start index of the resulting sub-vector in  $t_1$ 's partial output (i.e.,  $\{a, 0\}$ ). We only cache the left operand (sub-matrix DD node  $m_4$ ) because the right operand (*h*-sized sub-vector in *V*) is the same for all assigned multiplication tasks on the same thread. Similarly,  $t_2$  caches  $m_5$ and  $\{c, h\}$  (lines 8-10). At step ②,  $t_1$  and  $t_2$  execute  $bm_4 \cdot V[0:h]$ and  $dm_5 \cdot V[h:2h]$  simultaneously. Again, they check if  $m_4$  and  $m_5$  are cached, discovering historical results (lines 5-6).  $t_1$  finds an *h*-sized sub-vector in its partial output (stored in buffer B[0]) at start index 0 with weight *a*. It then multiplies this sub-vector by FlatDD: A High-Performance Quantum Circuit Simulator using Decision Diagram and Flat Array



Figure 7: DMAV with caching on four threads, illustrated in the four colors. The input matrix and vector are M and V, and the resulting vector is W (i.e.,  $M \cdot V \rightarrow W$ ). DD edges without labels have a weight of one by default.

b/a using SIMD-enabled scalar multiplication and places it in the fourth quarter of the B[0] (line 7). Likewise,  $t_2$  performs a similar scalar multiplication. Finally, at step (3), we sum the buffers to form the final result *W* using multi-threading and SIMD (lines 11-14).

3.2.3 Computational Cost in DMAV. Based on our observation, the MAC operation dominates DMAV computation. For example, within the Run function (Algorithm 1), which accounts for 99.99% of DMAV's runtime, the only non-trivial computation is performed by line 19, which executes a MAC operation. Therefore, we introduce the computational cost model based on the number of MAC operations.

First, it is necessary to determine the number of MAC operations in a DMAV. As shown in Figure 8, we use DFS to traverse DD nodes and a look-up table (MAC count table T) to track MAC operations for unique nodes, given that identical DD nodes incur the same number of MAC operations. The number of MAC operations of each node is the sum of that of its children, and the terminal node has one MAC operation. For each node, we traverse the outgoing edges from top to bottom and from left to right. Following this rule, at step (1), we go through  $m_1$ ,  $m_2$ ,  $m_3$ , and arrive at  $m_5$ .  $m_5$ only has one outgoing edge to a terminal node with one MAC (i.e.,  $T(m_5) = 1$ ). At step (2), we proceed to  $m_3$ 's lower-right outgoing edge, which also points to  $m_5$ . Thus,  $m_3$  has two MAC operations (i.e.,  $T(m_3) = 2T(m_5) = 2$ ). Then, we return to  $m_2$ , and visit the upper-right edge pointing to  $m_4$ . Similarly, we derive that  $T(m_6) =$ 1 and  $T(m_4) = 2$ , at steps (3) and (4). The lower outgoing edges of  $m_2$  connect to the previously visited nodes  $m_3$  and  $m_4$ , giving us  $T(m_2) = 2T(m_3) + 2T(m_4) = 8$  at step (5). Finally, at step (6), we return to  $m_1$ , obtaining  $T(m_1) = 2T(m_2) = 16$ . The total MAC count for this DMAV is  $T(m_1) = 16$ .

Then, we model the computational cost based on the number of MAC operations. We observe that DMAV's workload distribution among threads, with or without caching, is always balanced. Therefore, the computational cost model evenly divides the number of MAC operations by t threads. Suppose  $K_1$  is the number of MAC



Figure 8: Counting the number of MAC operations of a DMAV.

operations for a DMAV without caching, Equation 5 models its computational cost  $C_1$ .

$$C_1 = \frac{K_1}{t} \tag{5}$$

For DMAV with caching, we consider three costs. (1) **The cost** of scalar multiplication. By identifying *H* cache hits from repeated nodes in  $v_M$  (Algorithm 2), each corresponding to a scalar multiplication of size  $2^n/t$ , we determine that the number of MAC operations for scalar multiplication is  $2^nH/t$ . If SIMD computes *d* data elements simultaneously, on *t* threads, the cost of scalar multiplication is  $2^nH/(d \cdot t^2)$ . (2) **The cost of summing buffers.** Summing *b* buffers, each of size  $2^n$ , incurs  $2^n b$  MAC operations. When using SIMD on *t* threads, the cost of summing buffers is  $2^nb/(d \cdot t)$ . (3) **The cost unrelated to caching**. We count the MACs unrelated to caching in DMAV, as in Figure 8, this time avoiding the addition of the repeated nodes mentioned in (1). Suppose there are  $K_2$  MAC operations unrelated to caching, the corresponding cost is  $K_2/t$ . In summary, we model the computational cost  $C_2$ , for DMAV with caching as the following.

$$C_2 = \frac{K_2}{t} + \frac{2^n}{d \cdot t} (\frac{H}{t} + b)$$
(6)

To minimize the total computational cost, we choose DMAV with caching if  $C_1 > C_2$ , and without caching if not. To sum up, the computational cost for a DMAV operation is min $\{C_1, C_2\}$ .

#### 3.3 DMAV-Aware Gate-Fusion Algorithm

To further boost FlatDD's performance on large circuits, we propose a DMAV-aware gate-fusion algorithm that greedily fuses gates to reduce the total computational cost of DMAV.

Figure 9 shows the advantage of gate fusion. Consider two quantum gates  $M_1$ ,  $M_2$  in DD, there are two approaches to applying them consecutively to a state vector V. (1) **Sequential DMAV**: Multiply  $M_1$  by V to obtain intermediate state  $W_1$ , and then multiply  $M_2$  by  $W_1$  to obtain the final state  $W_2$ , as shown in Figure 9a. (2) **Gate fusion**: Apply DD-based matrix-matrix multiplication (DDMM) to  $M_1$  and  $M_2$ , fusing them into  $M_{21}$ , and then multiply  $M_{21}$  by V to obtain the final state  $W_2$ , as shown in Figure 9b. For simplicity, we evaluate two options using Equation 5 to determine which one has a lower computational cost. In sequential DMAV, the two consecutive DMAVs result in a total computational cost of 512/t. In contrast, gate fusion performs one DDMM and one DMAV. The DMAV computational cost is 256/t. The DDMM calls 12 multiplications and 4 additions, which is negligible compared to the DMAV

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Figure 10: Gate fusion can result in more computation.

computational cost since each multiplication or addition only constructs one DD node. Therefore, gate fusion reduces computation compared to sequential DMAV.

However, gate fusion does not always reduce computation, as shown in Figure 10. In Figure 10a, sequential DMAV without gate fusion incurs a computational cost of 6144/t. Conversely, in Figure 10b, gate fusion results in a DMAV computational cost of 8192/t and 21 DDMM multiplication calls. In this case, sequential DMAV has a lower computational cost.

To reduce the total computational cost, we fuse two gates only when the fused gate has a smaller computational cost compared to sequential DMAV. Accordingly, we introduce our DMAV-aware gate fusion in Algorithm 3. Algorithm 3 operates on the group of remaining gates, G, after FlatDD conversion. We initialize  $M_p$  (the previous gate matrix) to an identity DD matrix and set its DMAV computational cost (Section 3.2.3)  $C_p$  to 0 (line 2). Next, we iterate through G (line 3). For the *i*th gate matrix  $M_i$ , we calculate its DMAV computational cost,  $C_i$ . Then, we multiply  $M_i$  by  $M_p$  to form  $M_{ip}$ , and calculate its DMAV computational cost  $C_{ip}$  (line 4). If  $C_i + C_p <$  $C_{ip}$ , sequential DMAV is computationally cheaper than gate fusion. We then add  $M_p$  to the resulting gate group S. Before the next iteration, we assign the *i*th gate  $M_i$  and its DMAV computational cost  $C_i$  to the previous gate  $M_p$  and its DMAV computational cost  $C_p$  (line 6). If  $C_i + C_p < C_{ip}$  fails to satisfy, we fuse the current gate  $M_i$  with  $M_p$ . Then, we assign the fused gate  $M_{ip}$  and cost  $C_{ip}$  to the previous gate  $M_p$  and cost  $C_p$  (line 8). Finally, we return S at line 9.

#### 4 EXPERIMENTAL RESULTS

In this section, we evaluate the performance of FlatDD using 12 commonly used quantum circuits from QASMBench [69], MQT

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Algorithm 3 DMAV-aware gate fusion							
<b>Input:</b> <i>G</i> , group of remaining gates after FlatDD conversion.							
<b>Output:</b> <i>S</i> , group of gates after gate fusion.							
1: <b>function</b> GateFuse(G)							
2: $M_p \leftarrow I_{DD}; C_p \leftarrow 0$							
3: for $M_i \in G$							
4: $C_i \leftarrow \text{cost}(M_i); M_{ip} \leftarrow \text{DDMM}(M_i, M_p); C_{ip} \leftarrow \text{cost}(M_{ip})$							
5: <b>if</b> $C_i + C_p < C_{ip}$ <b>then</b>							
6: $S \cup \{M_p\}; C_p \leftarrow C_i; M_p \leftarrow M_i$							
7: else							
8: $M_p \leftarrow M_{ip}; C_p \leftarrow C_{ip}$							
9: return S							
10: end function							

Bench [88] and Quantum supremacy [7]. These benchmarks span both regular and irregular structures. First of all, we compare the runtime and memory performance of FlatDD with two state-ofthe-art QCSs (Section 4.2). Then, we demonstrate the scalability of FlatDD over increasing numbers of threads (Section 4.3). We then evaluate the effectiveness of our parallel DD-to-array conversion algorithm (Section 4.4) and DMAV caching technique (Section 4.5). Lastly, in Section 4.6, we evaluate our DMAV-aware gate-fusion algorithm on deep circuits with thousands of gates. All the experiments are conducted on a Ubuntu 22.04.2 LTS machine with 64 Intel Xeon Gold 6226R CPUs at 2.9 GHz and 256 GB memory capacity. We compile all programs with optimization flag -03 enabled. All SIMD operations are executed using Intel's AVX2 [60] vector instructions. For data with exponential difference, we measure the average in geometric mean.

## 4.1 Baselines

Given the large number of QCSs, it is not possible to compare FlatDD with all of them. Instead, we consider two representative QCSs, DDSIM [99] and Quantum++ [19], for two reasons: (1) Both DDSIM and Quantum++ are highly optimized and have demonstrated superior performance over existing QCSs. DDSIM introduces a DD-based compact representation of gate matrices and state vectors, as well as an efficient data structure for handling complex numbers [98]. On the other hand, Quantum++ leverages OpenMP [3] to enable multi-threaded simulation based on Eigen [23] array and matrix data structures. (2) Both DDSIM and Quantum++ are open-source [6, 87] and widely used by the community, allowing us to fairly study and reason their results.

#### 4.2 Overall Performance Comparison

Table 1 compares the overall performance of FlatDD with DDSIM and Quantum++. In this experiment, we do not incorporate the proposed gate-fusion algorithm but focus on the full-state simulation workload itself. We run FlatDD and Quantum++ using 16 threads and run DDSIM using one thread as DDSIM does not support multithreading. For all FlatDD runs, we set  $\beta = 0.9$  and  $\epsilon = 2$ , as these values are determined to be effective across multiple quantum circuits. Unless otherwise specified, all runtime and memory are measured in seconds (s) and megabytes (MB). We terminate the runs that take longer than 24 hours. We measure memory usage by recording the maximum resident set size (RSS) of our program using /bin/time. FlatDD: A High-Performance Quantum Circuit Simulator using Decision Diagram and Flat Array

On average, FlatDD is  $34.81 \times$  and  $17.31 \times$  faster over DDSIM and Quantum++ on all circuits. In terms of memory usage, FlatDD is  $1.70 \times$  and  $1.93 \times$  less than DDSIM and Quantum++. DDSIM demonstrates exceptional performance on regular circuits such as Adder and GHZ state. For instance, the circuit Adder has a very regular distribution of state amplitudes throughout the simulation; DDSIM takes less than 1 s while Quantum++ takes 1793.67 s. However, when the regularity does not appear frequently, such as DNN, VQE, and Quantum supremacy, DDSIM becomes significantly slower than FlatDD and Quantum++. For example, DDSIM is  $13.66 \times$  and  $12.64 \times$  slower than FlatDD and Quantum++ when simulating the 16-qubit DNN.

Although DDSIM and Quantum++ have their own strength in simulating certain circuits, FlatDD demonstrates a consistent performance advantage on all. For highly regular circuits, such as Adder, and GHZ state, where DDSIM performs extremely well, FlatDD also achieves fast runtime (< 1 s). This is because FlatDD does not switch from DDSIM to DMAV during the simulation. The only reason that FlatDD can be slower than DDSIM, such as simulating the 23-qubit GHZ state, is the overhead of DD size calculation and conversion timing checking (Section 3.1.1). For irregular circuits, such as 20-qubit DNN and quantum supremacy circuit, FlatDD can still take advantage of DD-based state vector to a certain extent before the state vector turns irregular. After this turning point, as observed from Figure 11a and Figure 11b, the runtime of DDSIM will significantly increase whereas FlatDD cleverly converts to DMAV and stays stable. Additionally, after the turning point, FlatDD is faster than Quantum++ primarily due to the efficient indexing pattern of our DMAV (see Section 3.2).

The memory usage of FlatDD is comparable to DDSIM and Quantum++. When the circuit is small, Quantum++ consumes the smallest memory compared with FlatDD and DDSIM, both of which require additional data storage for DDs. For example, when simulating the 16-qubit DNN, Quantum++ needs only 8.2 MB, whereas FlatDD and DDSIM need 32.26 MB and 69.6 MB, respectively. On the other hand, when the circuit is large, Quantum++ consumes significantly higher memory than FlatDD and DDSIM. For instance, when simulating the 25-qubit KNN, Quantum++ needs 1577.2 MB, while FlatDD and DDSIM only need 1078.78 MB and 388.1 MB, respectively. If the state vector is highly regular (e.g., Adder, GHZ state), FlatDD and DDSIM have similar memory usage, as FlatDD does not switch from DDSIM to DMAV. However, if the state vector contains high irregularity (e.g., DNN, VQE, Quantum supremacy), FlatDD has smaller memory usage than DDSIM because it will switch from DDSIM to DMAV to reduce DD overheads. Lastly, if the state vector is irregular in large circuits, FlatDD consumes smaller memory than both DDSIM and Quantum++. For example, when simulating the 26-qubit quantum supremacy circuit, FlatDD consumes 2132.48 MB memory, while DDSIM and Quantum++ consume 16799.4 and 3156.58 MB memory.

#### 4.3 Scalability of FlatDD

We demonstrate the scalability of FlatDD over increasing numbers of threads. Figure 12 illustrates the runtime of FlatDD and Quantum++ at different numbers of threads on two circuits (Quantum supremacy and KNN). When the number of threads increases,



Figure 11: Runtime comparison among FlatDD, DDSIM, and Quantum++ per gate.



Figure 12: Runtime scalability of FlatDD and Quantum++ under different numbers of threads.

FlatDD can complete the simulation faster, as illustrated in Figure 12a. For instance, at eight threads, FlatDD is 7.26× faster than one thread on KNN. FlatDD's performance saturates at about 16 threads. Likewise, in Figure 12b, Quantum++ shows a similar trend.

## 4.4 Evaluation of Parallel DD-to-Array Conversion Algorithm

Figure 13 compares the efficiency of FlatDD's parallel DD-to-array conversion algorithm with DDSIM on 10 quantum circuits using 16 threads. In this experiment, we integrate DDSIM's DD-to-array algorithm into FlatDD. Figure 13a compares the conversion time, and Figure 13b compares the cost of conversion in terms of its percentage in the total simulation time.

As shown in Figure 13a, FlatDD outperforms DDSIM in converting DDs to arrays on all circuits. For instance, our DD-to-array algorithm is 61.86× faster than DDSIM on KNN of 25 qubits; on average, FlatDD is 22.34× faster, which is attributed to multi-threading and SIMD parallelism. In terms of conversion cost, as shown in Figure 13b, FlatDD's DD-to-array algorithm takes only about 0.01–7% of the total runtime, while DDSIM can take up to 83.2% (Swap test of 25 qubits). This result highlights the efficiency of our parallel DD-to-array conversion algorithm.

#### 4.5 Evaluation of DMAV Caching Technique

Figure 14 demonstrates the efficiency of the proposed DMAV caching technique in terms of reduced computational cost and speed-up gain at different numbers of threads. We plot the results in a region (marked in blue) across the six largest quantum circuits,

Circuits			FlatDD (ours)			DDSIM [99]		Quantum++[19]			
Name	Qubits (n)	Gates	Runtime	Memory	Runtime	Speed-up	Memory	Runtime	Speed-up	Memory	
DNN	16	2032	4.12	32.26	56.27	13.66×	69.6	4.45	1.08×	8.2	
	20	6214	73.06	75.26	21847.81	299.02×	504.832	249.62	3.42×	61.34	
	25	9644	2283.02	1089.54	> 24 h	> 37.84×	≥ 4249.6	17527.8	7.68×	1586.54	
Adder	28	117	0.037	30.21	0.00167	0.045×	28.2	1793.67	$48477.57 \times$	12584.0	
GHZ state	23	46	0.028	30.21	0.004	0.14×	29.7	12.388	442.43×	397.93	
VQE	16	95	0.178	31.232	79.24	445.92×	50.2	0.20	1.12×	7.8	
KNN	25	39	4.75	1078.78	480.72	101.20×	388.1	135.669	28.56×	1577.2	
	31	48	500.22	67136.51	> 24 h	> 172.72×	≥ 2518.0	13123.5	26.23×	100657.37	
Swap test	25	39	4.84	1077.76	1649.81	340.94×	649.7	136.995	28.31×	1577.6	
Quantum supremacy	20	4500	101.38	65.54	67814.99	668.92×	700.928	185.207	1.83×	61.2	
	24	5560	1050.07	558.59	> 24 h	> 82.28×	≥ 5794.77	5034.7	4.80×	799.23	
	26	5990	3980.07	2132.48	> 24 h	> 21.71×	≥ 16799.4	22632.7	5.69×	3156.58	
Geor	Geometric mean		17.08	296.38	> 594.53	> 34.81×	≥ 504.52	295.62	17.31×	571.16	

Table 1: Comparison of simulation runtime and memory among FlatDD, DDSIM, and Quantum++ on 12 widely used circuits that exhibit different regularity structures.

Table 2: Comparison of simulation runtime and computational cost among FlatDD with DMAV-aware gate fusion (ours), FlatDD without gate fusion, and FlatDD with k-operations [100] on six deep circuits (> 1000 gates). Red. is the reduction in computational cost.

Circuits			FlatD DMAV-a fusion	D with ware gate (ours)	FlatDD without gate fusion				FlatDD with k-operations [100]			
Name	n	Gates	Runtime	Cost	Runtime	Speed-up	Cost	Red.	Runtime	Speed-up	Cost	Red.
DNN	16	2032	0.32	$5.4  imes 10^5$	4.12	12.81×	$8.2 \times 10^{6}$	15.4×	1.27	3.96×	$3.7 \times 10^{6}$	6.85×
	20	6214	5.09	$2.7 \times 10^{7}$	73.06	14.36×	$2.4 \times 10^{8}$	8.56×	25.858	5.1×	$1.1 \times 10^{8}$	3.89×
	25	9644	126.48	$1.2 \times 10^{9}$	2283.02	18.0×	$1.7 \times 10^{10}$	13.4×	688.396	5.44×	$5.3 \times 10^{9}$	4.32×
Quantum supremacy	20	4500	7.68	$4.7  imes 10^7$	101.38	13.2×	$3.8 \times 10^{8}$	8.04×	36.99	4.82×	$2.9 \times 10^{8}$	6.19×
	24	5560	90.90	$9.0 \times 10^{8}$	1050.07	11.55×	$7.4 \times 10^{9}$	8.26×	581.64	6.40×	$5.7 \times 10^{9}$	6.37×
	26	5990	405.04	$3.9 \times 10^{9}$	3980.07	9.83×	$3.2 \times 10^{10}$	8.21×	2568.12	6.34×	$2.6 \times 10^{10}$	6.73×
Geometric mean		19.70	$1.2 \times 10^8$	257.45	13.1×	$1.2 \times 10^{9}$	9.94×	103.83	5.27×	$6.7 \times 10^{8}$	5.59×	

DNN (n = 16, 20, 25) and quantum supremacy (n = 20, 24, 26), with a line showing the average. As shown in Figure 14a and Figure 14b, our caching technique can reduce the computational cost as the number of threads increases. For instance, with 16 threads, we can achieve 13.53% reduction. A similar trend can be observed in the speed-up plot (Figure 14b). For example, our caching technique contributes to an average of 16.47% speed-up at 16 threads where the performance saturates (see Figure 12). This result highlights the effectiveness of our caching technique under different numbers of threads.

## 4.6 Evaluation of DMAV-Aware Gate Fusion

In this experiment, we evaluate the performance advantage of our gate-fusion algorithm by running FlatDD on the six deepest circuits (DNN and supremacy) with and without gate fusion. We compare the result with k-operations [100], a particularly effective gate-fusion algorithm designed for DD-based simulators. As shown in Table 2, our gate-fusion algorithm achieves 13.1× and 5.27× speedup in terms of geometric mean when compared with FlatDD without gate fusion and k-operations, respectively. Our gate-fusion algorithm also reduces computational cost by  $9.94\times$  and  $5.59\times$  when compared with FlatDD without gate fusion and k-operations, respectively. This is because our gate-fusion algorithm always fuses a quantum gate that reduces the computational cost. The result in

Table 2 shows the effectiveness of our DMAV-aware gate-fusion algorithm.

#### 5 CONCLUSION

In this paper, we have introduced FlatDD, a parallel quantum circuit simulator that combines the strengths of DD-based and array-based simulators. Evaluated on commonly used quantum circuits, FlatDD has demonstrated 34.81× speed-up and 1.93× memory reduction compared to state-of-the-art simulators. Inspired by our success of GPU-accelerated computing [12–18, 24–33, 38–58, 61, 62, 64–67, 70–80, 83, 84, 92, 93, 96], our future work will enhance the simulation performance using GPU.

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(b) Conversion percentage in total runtime comparison.

# Figure 13: Comparison between FlatDD's parallel DD-to-array algorithm and DDSIM's DD-to-array algorithm.



Figure 14: Performance comparison between DMAV with caching and DMAV without caching over different numbers of threads on various quantum circuits.

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