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A Hardware-Aware Heuristic for the Qubit Mapping Problem in the NISQ Era

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ABSTRACT Due to several physical limitations in the realization of quantum hardware, today's quantum computers are qualified as noisy intermediate-scale quantum (NISQ) hardware. NISQ hardware is characterized by a small number of qubits (50 to a few hundred) and noisy operations. Moreover, current realizations of superconducting quantum chips do not have the ideal all-to-all connectivity between qubits but rather at most a nearest-neighbor connectivity. All these hardware restrictions add supplementary low-level requirements. They need to be addressed before submitting the quantum circuit to an actual chip. Satisfying these requirements is a tedious task for the programmer. Instead, the task of adapting the quantum circuit to a given hardware is left to the compiler. In this article, we propose a hardware-aware (HA) mapping transition algorithm that takes the calibration data into account with the aim to improve the overall fidelity of the circuit. Evaluation results on IBM quantum hardware show that our HA approach can outperform the state of the art, both in terms of the number of additional gates and circuit fidelity.

INDEX TERMS Noisy intermediate-scale quantum (NISQ) hardware, quantum computing, qubit mapping.

NOMENCLATURE

Notation *Definition*

q	Logical qubits for quantum circuit.
Q	Physical qubits for quantum device.
g	Quantum gate.
$g.q_n$	n th logical qubit the quantum gate g is applied on.
G	Hardware coupling graph.
D	Distance matrix.
S	Swap matrix.
$G_{\mathcal{E}}$	Hardware graph with swap error rates as weights.
\mathcal{E}	Swap error matrix.
G_T	Hardware graph with swap execution time as weights.
T	Swap execution time matrix.
H	Heuristic cost function.
π	Mapping from q to Q .
F	First layer.
E	Extended layer.
W	Weight parameter.

I. INTRODUCTION

In recent years, quantum computing has become a very active field of research. It promises to solve classically intractable computational problems such as integer factorization [1], quantum chemistry [2], linear algebra [3]–[8], or optimization [9]–[11]. Along with algorithms, quantum hardware has attracted the attention of several companies such as IBM, Google, Intel, or Rigetti that have demonstrated quantum chips with 53, 72, 49, and 28 qubits, respectively. IBM and Rigetti have also given access to a cloud quantum computing service on which anyone can submit quantum circuits to real quantum hardware. The aforementioned quantum hardware can already be qualified as noisy intermediate-scale quantum (NISQ) hardware [12]. Still, none of them is fault-tolerant as quantum error correction codes are in infancy. Nevertheless, it is believed that even a noisy quantum chip with limited qubit-to-qubit connectivity can be used to solve some classically intractable problems, one of the most promising candidates being quantum chemistry [2].

From the algorithm perspective, a new paradigm for quantum algorithms has emerged to take into account the limitations of NISQ hardware—variational algorithms. Examples of variational algorithms include the variational quantum eigensolver (VQE) [13], the variational quantum linear solver (VQLS) [7], [8], or the quantum approximate optimization algorithm (QAOA) [11]. However, there is a difference between the quantum program written by the programmer and what can be executed on the current quantum hardware. Quantum programs are written as if they were running on ideal quantum hardware without any noise or physical constraints. But real quantum chips are not ideal—for example, for superconducting devices which are targeted in this article, current two-qubit gates can at best only be applied between two neighboring qubits. If we want to perform quantum computations, our quantum circuits must obey such connectivity constraints, which mean that a modification of the quantum program is necessary to adapt it to the real quantum device. This problem of adapting a quantum program to given hardware connectivity is called the qubit mapping problem and is the focus of this article.

The qubit mapping problem can be reformulated as two subproblems. First, to find an initial mapping, i.e., a mapping between the “logical qubits” (as a qubit in a quantum circuit) to the “physical qubits” (as a qubit in a quantum chip). Second, to determine a mapping transition algorithm to identify the quantum gates to insert in a quantum circuit such that it complies with the targeted quantum hardware topology. Finding the optimal solution for the qubit mapping problem is likely to be an NP-complete problem as noted in [14].

Two types of methods have been used to solve the qubit mapping problem. The first method is to reformulate it as a mathematically equivalent problem that can then be solved using a specialized solver. Such mathematical formalism can be integer linear programming [15]–[18], satisfiability modulo theory [19], [20], or even constraint programming [21], [22]. However, these mathematical approaches suffer from long runtime and are difficult to scale up. The second method is to use heuristics to modify the quantum circuit, starting from the first quantum gate and transforming the circuit sequentially by making each gate one after the other hardware-compliant.

Most of the previous works [23]–[27] using the second method only adapt for nearest-neighbor connectivity and cannot be directly applicable to actual quantum architectures with nonuniform connections. Recently, publications [14], [28]–[34] that are not restricted to a specific architecture have been released. The algorithm presented in [28] uses a heuristic to find the best permutation at each step of the mapping procedure. Instead of representing a quantum circuit as a fixed sequence of layers, [35] introduces the directed acyclic graph (DAG) that takes into account the dependency and commutativity of quantum gates. A major improvement has been shown by [29], which uses a “forward–backward–forward” mapping algorithm. Moreover, the “lookahead” strategy has been introduced in the heuristic cost function for

further optimization in some existing works, notably [28]–[31], [34]. For qubit movement, most of these methods only use SWAP gate. A notable exception is [31] that considered both Bridge and SWAP gate. Moreover, most of these works aim to minimize the number of inserted gates and do not consider the noise impact on different qubits.

In [16], [19], [36]–[38], calibration data are exploited and the applied approach is to insert additional gates between strongly linked qubits, i.e., qubits linked with a low two-qubit gate error rate. However, these works do not consider a holistic view of the problem such as exploring initial mapping or the heuristic function is not efficient enough to select the best candidate of the inserted gate.

In this article, we follow the second type of methods that consist in developing a heuristic to choose the best SWAP to insert based on calibration data. We propose a hardware-aware (HA) heuristic mapping transition algorithm to address the drawbacks mentioned above. Our main contributions can be listed as follows. First, we present a mapping transition algorithm that takes into account the hardware topology and the calibration data to improve the overall output state fidelity and reduce the total execution time. Second, to reduce the number of additional gates required to map the quantum circuit to the quantum chip, our algorithm can select between a SWAP or Bridge gate. Finally, we run our HA algorithm on real quantum hardware and compare with various mapping methods from the literature.

II. STATE OF THE ART

Here, we introduce state of the art on quantum hardware devices and their constraints, focusing mainly on IBM quantum devices. Then, we explain the qubit mapping problem. Finally, a small motivational example is shown to illustrate the gist of our algorithm. The notations used in this article are summarized in the nomenclature (we reference some notations from [29]).

A. CURRENT STATE OF THE ART ON QUANTUM HARDWARE

NISQ hardware is characterized in [12] as quantum hardware having from 50 to a few hundred noisy qubits on which one can only perform noisy operations. At the time of writing, several companies have already demonstrated quantum chips that can, according to the definition, be qualified as NISQ chips. For example, IBM announced the latest 53-qubit quantum chip and gave access to the community to execute quantum circuits. Other companies like Google (72 qubits) or Intel (49 qubits) announced quantum chips that could be qualified as NISQ but did not provide any information about their characteristics. One of the significant challenges of these quantum chips that limits them from solving real-world problems is their level of noise—even if these chips have enough qubits theoretically to show a quantum speedup, the fidelity of their quantum operations is still too low to obtain any advantage over the classical computer on real-world

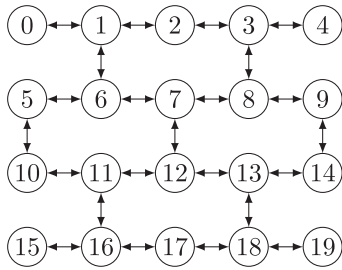


FIGURE 1. `ibmq_almaden` topology. Qubits are represented as circles and indexed from 0 to 19. A connection between two qubits is represented by an edge between the two qubits.

TABLE 1. `ibmq_almaden` characteristics

Qubit number	20
Single qubit error rate	$2.655e^{-4}$ to $9.569e^{-4}$
CNOT error rate	$8.136e^{-3}$ to $3.403e^{-2}$
id gate length	35.56 ns
u1 gate length	0 ns
u2 gate length	35.56 ns
u3 gate length	71.11 ns
CNOT gate length	248.88 ns to 860.44 ns
T1	34.66 μ s to 139.46 μ s
T2	12.16 μ s to 200.25 μ s

Note that the exact hardware characteristics are not constant and change at each recalibration of the chip.

problems. In this article, we mainly focus on IBM architectures. Other hardware such as Google's Sycamore [39] or Rigetti's Aspen-7 is not specially targeted. Still, the proposed algorithm and methods are general enough to be applicable to any quantum chip that use the quantum-gate model of computation and so should be applicable to this hardware.

Fig. 1 shows the topology, also called coupling graph, of IBM Quantum's `ibmq_almaden`, a 20-qubit system. Each vertex represents a qubit and the edge represents the coupling interconnect between two qubits. Table 1 shows the calibration data that are extracted from [40]. It includes CNOT error rates, single qubit error rates, energy relaxation and decoherence characteristic times T1 and T2, and execution time (gate length). The calibration data show that the error of two-qubit gates is one order of magnitude higher than their one-qubit counterparts. This is also the case for gate execution times—two-qubit gates are approximately an order of magnitude slower than one-qubit gates. For simplicity and because of the relatively low error rates and execution times of one-qubit gates when compared to two-qubit gates, we focus on two-qubit gates in this article.

Moreover, it is important to note that all the interconnects between qubits are not equal with respect to CNOT gate error rate or execution time. Taking `ibmq_almaden` as an example, the best CNOT gate has an error rate of 4.18 times lower than the worst CNOT and the maximum execution time is

3.46 times longer than the minimum one. Therefore, we cannot treat each qubit equally, and we need to consider the interconnect topology between qubits as well as their error rate. CNOT gates can be applied in either direction by conjugating with H gates. As we do not consider one-qubit gates in this study, we do not have to consider the connectivity direction.

B. QUBIT MAPPING PROBLEM

Following the abstraction first introduced in classical computing decades ago, most of the quantum circuits are described in a generic manner that does not take into account all the physical hardware constraints. Many of the currently existing frameworks for quantum algorithm development encourage this way of development by giving access to a broad set of “primitive” gates. For example, the Qiskit library allows the developer to choose from more than 30 primitive gates, whereas the IBM quantum chips only provide four physical hardware gates (five if we take the identity gate into account). However, any gate can also be implemented with OpenPulse [41], which is a low level hardware control for users to generate their gates to mitigate errors. Such an abstraction relieves the burden from the developer to adapt the code to a specific hardware and transfer it to the compiler, whose role is to transform an abstracted code into the most efficient hardware code possible. To do so, a compiler for quantum programs should perform several steps summarized in the following paragraphs.

The first step is to decompose the abstracted quantum gates into hardware gates. The hardware gates available strongly depend on the quantum hardware we are compiling for, but are generally comprised of a two-qubit “entangler” gate (controlled-X gate for IBM hardware, fSim gate for Google hardware) and several one-qubit gates (u1, u2, u3, and id gates for IBM hardware). At the end of this step, the quantum circuit has been modified to only contain quantum gates that are directly implemented in the quantum hardware.

But translating all abstract gates to hardware gates is generally not enough to make the quantum circuit executable on the specific hardware—the hardware topology is rarely respected at the end of this first step and the circuit requires a second step with further modifications. Such modification of the quantum circuit to make it compliant with the hardware topology is often done by inserting SWAP gates before nonexecutable two-qubit gates. Note that on current hardware, only two-qubit gates are restricted by the hardware topology.

Finally, once the quantum circuit is executable on the specified quantum hardware, a final third step is performed to optimize the quantum circuit. Depending on the figure of interest, the optimization can aim at reducing the execution time, gate count, increasing the final state fidelity or even reducing the number of qubits needed.

The qubit mapping problem is defined as the second compilation step that modifies the quantum circuit to contain only two-qubit gates that fit into the hardware topology. But in practice qubit mapping algorithms also try to consider the

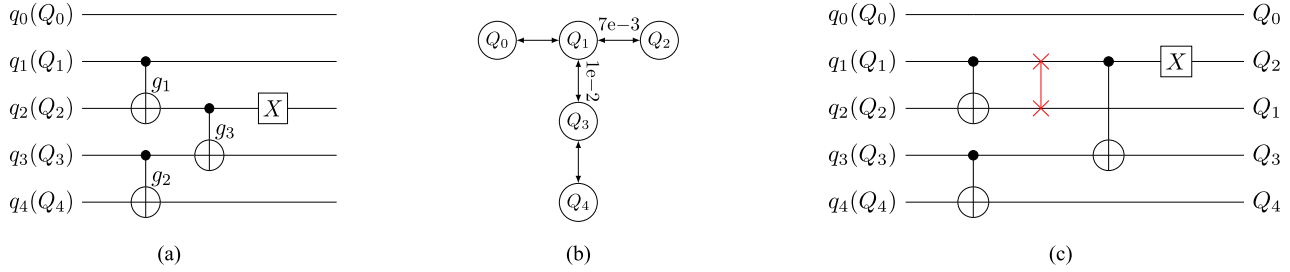


FIGURE 2. Motivational example for the qubit mapping problem. (a) Original circuit. (b) *ibmq_valencia*. (c) Updated circuit.

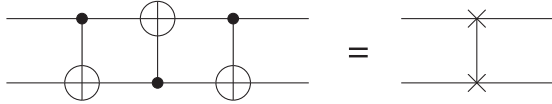


FIGURE 3. SWAP gate.

third step that consists in optimizing the generated quantum circuit according to a chosen figure of merit.

C. MOTIVATIONAL EXAMPLE

Fig. 2(a) shows a small quantum circuit, which is composed of three CNOT gates and one X gate. It is mapped to a 5-qubit IBM quantum device called *ibmq_valencia*, shown in Fig. 2(b). For simplicity, the initial mapping is allocated linearly as $\{q_0 \rightarrow Q_0, q_1 \rightarrow Q_1, q_2 \rightarrow Q_2, q_3 \rightarrow Q_3, q_4 \rightarrow Q_4\}$. Gates g_1 and g_2 comply with the hardware topology (i.e., coupling constraints) and can be executed directly. However, g_3 is applied to two nonconnected qubits. Therefore, a movement (i.e., a SWAP gate, shown in Fig. 3) of logical qubits is needed before being able to execute g_3 on the hardware connection between q_2 and q_3 . Referring to the coupling graph in Fig. 2(b), three SWAP gates are possible: $\{q_1, q_2\}$, $\{q_1, q_3\}$, and $\{q_3, q_4\}$. Among these possible SWAPs, two of them change the current mapping between logical and physical qubits in such a way that the CNOT gate between q_2 and q_3 becomes executable—swapping of $\{q_1, q_2\}$ and $\{q_1, q_3\}$. Translating the logical qubits to their physical counterparts, the SWAPs $\{Q_1, Q_2\}$ and $\{Q_1, Q_3\}$ are our candidates. At this step, most of the state-of-the-art algorithms consider the two possible SWAPs to be equal and will randomly select one. However, if the calibration data are considered, the SWAP between $\{Q_1, Q_2\}$ is less noisy than the other [error rate of the two interconnects is shown in Fig. 2(b)]. A SWAP operation consists of three CNOTs and we want to insert a SWAP gate with the least noise. Thus, the SWAP gate between $\{q_1, q_2\}$ is inserted and the final mapping is $\{q_0 \rightarrow Q_0, q_1 \rightarrow Q_2, q_2 \rightarrow Q_1, q_3 \rightarrow Q_3, q_4 \rightarrow Q_4\}$. The updated circuit is shown in Fig. 2(c).

III. PROPOSED SOLUTION

We are inspired by the SABRE algorithm presented in [29], which is a SWAP-based heuristic algorithm to reduce the number of additional CNOT gates. We propose a HA SWAP

and Bridge based heuristic search algorithm. Compared to SABRE algorithm, which aims at reducing the number of additional gates, we improve the circuit fidelity as well as reduce the number of additional gates by introducing a new distance matrix that takes into account both of the hardware connectivity and the calibration data. Moreover, SABRE only uses SWAP gate when a qubit movement is needed, whereas our algorithm decides between a SWAP and Bridge for qubit movement to further reduce the number of additional gates. Finally, we also develop an initial mapping algorithm called hardware-aware simulated annealing (HSA) in order to evaluate the mapping transition algorithm of different flavors.

The compiler takes as input a quantum program written in the OpenQASM language [42] and the calibration data of a specific IBM quantum device. During the compilation process, it considers the hardware constraints such as hardware topology and gate availability. Then, the qubit mapping algorithm is applied. It contains two principal parts—initial mapping and mapping transition algorithm. In the mapping transition step, some optimizations are done to generate a circuit with a better performance in terms of final state fidelity. The source code is publicly available at <https://github.com/peachnuts/HA>.

We start by explaining our HA algorithm in Section III-A. In Section III-B, we describe the HSA method for initial mapping. Finally, Section III-C presents the metrics used to evaluate our algorithm.

A. HA SWAP AND BRIDGE-BASED HEURISTIC SEARCH

The first step of the algorithm is to process the input quantum circuit in order to reformulate it in a more convenient data format. Starting from the input quantum circuit, we can obtain a DAG circuit, which represents the operation dependencies in the quantum circuit without considering the hardware constraints. The DAG is constructed such that quantum gates are represented by the graph nodes and the directed edge (i, j) between nodes i and j represents a dependency from gate i to j , i.e., gate i should be executed before j .

Once the DAG is constructed, graph nodes (i.e., quantum gates) can be ordered according to the gate dependencies—for example, if gate j depends on gate i , then gate i will be ordered before gate j . One possible ordering that fulfill this property of dependency is the well-known topological

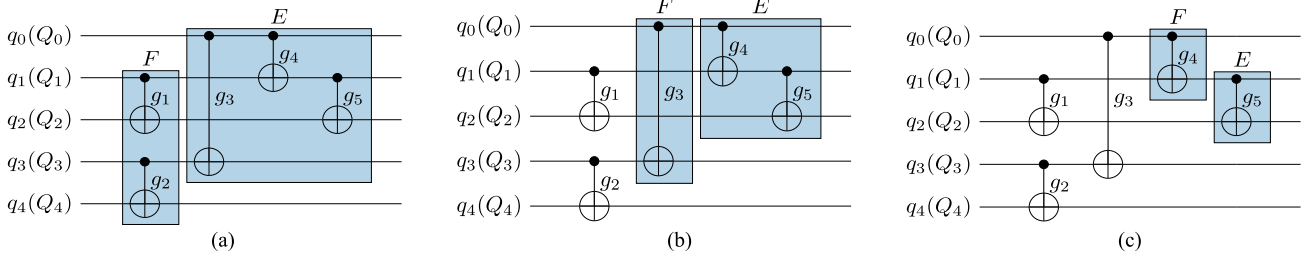


FIGURE 4. Evolution of the layers F and E on a simple circuit with a detailed explanation at each step. (a) Beginning of the HA mapping algorithm. g_1 and g_2 do not overlap and are the first gates in the circuit so they are in F . The other gates are pushed in E . (b) g_1 and g_2 are compliant with the hardware topology. They are executed and removed from F . The gate g_3 is pushed into F but is not compliant and a SWAP/Bridge should be inserted. g_4 overlaps with g_3 and cannot be inserted in F . (c) After Bridge insertion, g_3 is executed and removed from F . g_4 no longer overlap with a gate in F and is added to the first layer. g_5 overlaps with g_4 and so should stay in E .

ordering. Note that depending on the quantum circuit, this ordering might not be unique.

Quantum gates can then be divided into three groups: 1) the executed gates; 2) the executable gates; and 3) to-be-executed future gates. Executed gates are quantum gates that have already been mapped by the algorithm. Executable gates constitute the first layer, denoted F . A gate is considered executable when all the gates it depended on are in the executed gates group. Finally, to be executed future gates are the rest of the gates (not yet executed nor executable). These gates are included in the extended layer, E . An illustration of layers E and F is shown in Fig. 4.

1) HEURISTIC COST FUNCTION

A heuristic cost function H is introduced to estimate the cost of each possible (i.e., executable) swap pairs at a given step of the iterative algorithm. Its objective is to quantify the quality of the possible swap pairs according to the distance considered and to select the best swap pair.

When inserting a SWAP gate, the circuit is divided into two layers: 1) the first layer F ; and 2) the extended layer E . Note that inserting a SWAP gate will not only influence the gates in the first layer F but also the gates in the extended layer E . The approach of considering the swap pair's impact on the extended layer is referred as the lookahead ability. It can contribute to a better selection and depends on the size of the extended layer.

We devise several metrics that can be used to estimate the cost of a swap pair in HA. We consider three different distance matrices—swap matrix S , swap error matrix \mathcal{E} , and swap execution time matrix T . Because S , \mathcal{E} , and T contain entries with incompatible units and different scales, we update T to make it dimensionless and each matrix is normalized. Moreover, we introduce weights (α_1 , α_2 , and α_3 for S , \mathcal{E} , and T , respectively) to allow to choose the importance of each parameter in terms of number of SWAPs, gate error and execution time.

Matrix S is constructed such that the entry (i, j) stores the distance on the real hardware between qubit i to a neighbor of qubit j , which is also equal to the minimum number of SWAP gates needed to move qubit i to qubit j . The matrix

is efficiently constructed by using the Floyd–Warshall algorithm [43].

Matrix \mathcal{E} stores in its entry (i, j) the minimum error rate attainable to move the qubit i to a neighbor of qubit j . The error rate of each possible SWAP is computed based on the calibration data of CNOT gates. The decomposition of a SWAP gate in terms of CNOT gates is shown in Fig. 3.

The success rate of a CNOT between the physical qubits Q_i and Q_j , denoted by $S(Q_i, Q_j)$, is computed from the error rates given in the calibration data. Equation (1) computes the error rate of a SWAP gate between two connected physical qubits Q_i and Q_j while taking into account that the swap operation is symmetric. The final \mathcal{E} matrix is constructed by using the Floyd–Warshall algorithm on the graph $G_{\mathcal{E}}$ with the computed errors as edge weights.

$$G_{\mathcal{E}}(Q_i, Q_j) = 1 - S(Q_i, Q_j) \times S(Q_j, Q_i) \times \max(S(Q_i, Q_j), S(Q_j, Q_i)) \quad (1)$$

Matrix T is computed, similarly as S and \mathcal{E} , with the Floyd–Warshall algorithm applied on graph G_T but by using the SWAP execution time. This execution time is computed with (2), where $t(Q_i, Q_j)$ is the execution time of the CNOT gate with Q_i as control and Q_j as target, extracted from the calibration data.

$$G_T(Q_i, Q_j) = t(Q_i, Q_j) + t(Q_j, Q_i) + \min(t(Q_i, Q_j), t(Q_j, Q_i)) \quad (2)$$

The summation of the three matrices forms a new matrix called distance matrix D [shown in (3)]. The distance matrix represents the “distance” between each pair of qubits in the quantum chip. Here, the “distance” means the combination of swap distance, overall error rate, and execution time of the shortest path.

$$D = \alpha_1 \times S + \alpha_2 \times \mathcal{E} + \alpha_3 \times T \quad (3)$$

Inserting a SWAP gate will have an impact on the current mapping π_c , changing it to π_{temp} . We compute the cost of this SWAP on the first layer F with the cost function H_{basic} shown in (4). A small score means the SWAP has a little impact on the first layer gates with respect to the overall

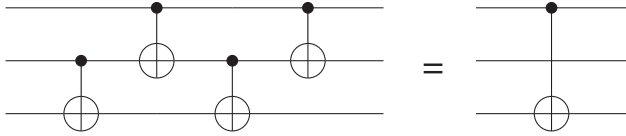


FIGURE 5. Bridge gate.

distance considered. The swap pair with the minimum score is selected as the best candidate.

$$H_{\text{basic}} = \sum_{g \in F} D[\pi_{\text{temp}}(g.q_1)][\pi_{\text{temp}}(g.q_2)] \quad (4)$$

We also consider the impact of the swap pair on the extended layer E . The impact of a SWAP on the first layer is prioritized over its impact on the extended layer. As a result, a weight parameter W is added to the extended layer cost to scale its impact. Moreover, the impacts on the first layer and extended layer are normalized by dividing them with their respective number of gates. The complete heuristic function including the extended layer E with lookahead ability is shown in (5). Even though (4) and (5) are similar to equations in [29], it is important to note that the distance matrix D is different.

$$H = \frac{1}{|F|} \sum_{g \in F} D[\pi_{\text{temp}}(g.q_1)][\pi_{\text{temp}}(g.q_2)] + W \times \frac{1}{|E|} \sum_{g \in E} D[\pi_{\text{temp}}(g.q_1)][\pi_{\text{temp}}(g.q_2)] \quad (5)$$

2) SWAP GATE AND BRIDGE GATE

Another important metric of the HA algorithm is the heuristic cost function that estimates the usefulness of a SWAP. In some situations, even the best SWAP may have a negative impact on the overall circuit. In that case, inspired by [31], our heuristic function decides to insert a Bridge gate instead of a SWAP gate if the topology allows it. The decomposition of the Bridge gate with four CNOTs is shown in Fig. 5. The Bridge gate allows executing a CNOT between two qubits that share a common neighbor. Both SWAP and Bridge gate need three supplementary CNOTs. Note that the Bridge gate can only be used to replace a CNOT if the distance between the control and target qubits (i.e., the minimum number of links between the two qubits) is exactly two.

Fig. 6(a) shows an example of quantum circuit that is mapped to `ibmq_valencia` with the topology described in Fig. 2(b). The quantum gates g_1 and g_2 comply with the topology of the chip, but g_3 does not. By evaluating the heuristic cost function H , the SWAP between q_0 and q_1 is selected. But as shown in Fig. 6(b), the chosen SWAP has a negative impact on the extended layer—gate g_5 is no longer executable and another SWAP gate is required to execute it.

Such situations can be solved by using a Bridge gate instead of a SWAP gate as shown in Fig. 6(c). Since the

distance between the control qubit q_0 and the target qubit q_3 of gate g_3 is two, we can insert a Bridge gate instead. Using a Bridge gate allows to execute the CNOT gate g_5 without changing the current mapping. Moreover, by using a Bridge gate, we only add three CNOTs to map the entire circuit, instead of six (two times more) if only SWAP gates were used.

Once the cost H of each swap pair is computed, the heuristic will try to choose the best option between inserting a SWAP or Bridge gate. To do so, it considers two mappings: π_c , the mapping used before selecting the best swap pair and also the mapping obtained after inserting a Bridge gate, and π_{temp} , the new mapping that would be obtained after inserting the best SWAP gate. The overall effect of the SWAP gate on the extended layer E is computed according to (6). If the effect of the best SWAP gate is negative, this means that the considered swap pair has an overall negative impact on the extended layer E . In this case, we consider that it is better to keep the current mapping so, if the hardware topology permits it, a Bridge gate is inserted instead of a SWAP gate.

$$\text{Effect} = \sum_{g \in E} D[\pi_c(g.q_1)][\pi_c(g.q_2)] - D[\pi_{\text{temp}}(g.q_1)][\pi_{\text{temp}}(g.q_2)] \quad (6)$$

3) HA ALGORITHM

The mapping transition algorithm will go through each quantum gate sequentially and mark the directly executable gates as executed. If no more gates can be marked as *executed*, this means that either the quantum circuit is fully mapped or all the gates in the first layer do not comply with the hardware topology. In the first case, the mapping algorithm can be stopped and the mapped quantum circuit returned. In the second case, the algorithm calls a heuristic function to choose the best SWAP or Bridge gate to insert in order to make some of the gates in the first layer executable. The algorithm then iterates, until the quantum circuit is fully mapped.

Algorithm 1 shows the pseudocode of the HA heuristic method. Note that the most recent calibration data should be retrieved (i.e., through the IBM quantum experience) before each usage of the HA algorithm to ensure that the algorithm has access to the most accurate and up-to-date information possible.

The heuristic method to insert a SWAP or Bridge when no gate in the first layer F is executable can be described as follows. First, a list of all the candidate SWAP gates, `swap_candidate_list`, is constructed based on the quantum gates in the first layer F and the hardware coupling graph G . Then, for each SWAP candidate a temporary mapping π_{temp} is computed with the `Map_Update` function. The final cost of the candidate SWAP is computed following (5). The SWAP with the minimum score is selected and called `swap_min`.

The last step is to choose between a SWAP gate or a Bridge gate. A SWAP gate can always be used, whereas

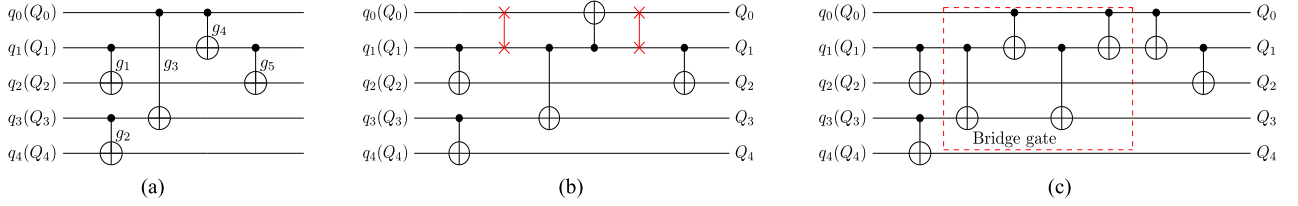


FIGURE 6. Example of a quantum circuit showing the difference between SWAP and Bridge transformation. (a) Original circuit. (b) SWAP gate transformation. (c) Bridge gate transformation.

a Bridge gate can only be inserted if a gate in the first layer F becomes executable from the mapping obtained after applying the swap_{\min} gate. If a Bridge gate is not insertable, then the algorithm has no choice but to insert a SWAP gate. Else, the algorithm decides the gate (SWAP or Bridge) to insert based on the effect of the SWAP gate on the extended layer computed with (6). If adding a SWAP gate has a negative impact on the extended layer, then a Bridge gate (which does not change the current mapping) is inserted. Otherwise, if adding a SWAP gate has a positive effect on the extended layer, then the algorithm inserts a SWAP gate.

4) RUNTIME ANALYSIS

The HA algorithm outperforms SABRE algorithm thanks to several modifications while not changing its asymptotic complexity. The mapping procedure is separated into two steps: an initialization step that is independent of the mapped quantum circuit and a mapping step.

The initialization step computes the distance matrix that is used afterward in the mapping step. In our algorithm, the distance matrix is computed according to (4). Each of S , \mathcal{E} , and T constituting the distance matrix D requires to use the Floyd–Warshall algorithm once on the hardware graph. This means that we need to perform three calls to an algorithm of $O(n^3)$ complexity, n being the number of qubits of the targeted quantum chip. Moreover, the weights used by the Floyd–Warshall algorithm for the matrices \mathcal{E} and T should be retrieved online with Qiskit API. This retrieval is an operation that theoretically takes $O(n^2)$ time in the worst case as we need to retrieve CNOT error rates and execution time for each link. Note that the current quantum chips only have $O(n)$ links and so the asymptotic complexity of this step is $O(n)$. Overall, the initialization step is dominated by the cost of applying the Floyd–Warshall algorithm, that takes $O(n^3)$ time.

After the initialization step, the actual mapping procedure is applied. Let n be the number of qubits, g the number of CNOT gates in the mapped quantum circuit and d the diameter of the chip, i.e., the minimum SWAP distance between the two farthest qubits on the quantum chip. In the worst case, all the CNOT gates should be mapped because none of them comply with the hardware topology. Moreover, all the CNOT gates might need up to d SWAPs in order to become executable. Finally, for each SWAP insertion, we need to execute the heuristic cost function. This function will need to

Algorithm 1: Heuristic Algorithm for Selecting Additional Gate Candidate.

input : Circuit DAG , Coupling graph G , Current mapping π_c , Distance matrix D , Swap matrix S , First layer F , Extended layer E , Weight parameter W

output: New mapping π_n , Inserted gate g_{add}

```

1 begin
2   Set score to empty list;
3   Set effect to empty list;
4   swap_candidate_list  $\leftarrow$  FindSwapPairs( $F$ ,  $G$ );
5   for swap  $\in$  swap_candidate_list do
6      $\pi_{\text{temp}} \leftarrow \text{Map\_Update}(\text{swap})$ ;
7      $H_{\text{basic}} \leftarrow 0$ ;
8     for gate  $\in F$  do
9        $H_{\text{basic}} \leftarrow H_{\text{basic}} + D(\text{gate}, \pi_{\text{temp}})$ ;
10    end
11     $H_{\text{extended}} \leftarrow 0$ ;
12    for gate  $\in E$  do
13       $H_{\text{extended}} \leftarrow H_{\text{extended}} + D(\text{gate}, \pi_{\text{temp}})$ ;
14      effect_cost  $\leftarrow$  effect_cost +  $D(\text{gate}, \pi_c) - D(\text{gate}, \pi_{\text{temp}})$ ;
15    end
16     $H \leftarrow \frac{1}{|F|} H_{\text{basic}} + \frac{W}{|E|} H_{\text{extended}}$ ;
17    score.append( $H$ );
18    effect.append(effect_cost);
19  end
20  Find the swap with minimum score:  $\text{swap}_{\min}$ ;
21  Find the gate in  $F$  that become executable by
    applying  $\text{swap}_{\min}$ :  $g_s$ ;
22  if effect[ $\text{swap}_{\min}$ ]  $< 0$  and  $S(g_s, \pi_c) = 2$  then
23     $\pi_n \leftarrow \pi_c$ ;
24     $g_{\text{add}} \leftarrow g_B$ ;
25  else
26     $\pi_n \leftarrow \text{Map\_Update}(\text{swap}_{\min})$ ;
27     $g_{\text{add}} \leftarrow \text{swap}_{\min}$ ;
28  end
29  return  $\pi_n, g_{\text{add}}$ ;
30 end

```

explore at most n^2 links (in the case of an all-to-all connected chip, this number improves to $O(n)$ on practical quantum chips with a nearest-neighbor connectivity), where exploring one link might take a time of $O(g)$ if all the CNOT gates are

included in either F or E . In summary, the mapping step takes $O(g^2 dn^2)$ time in the worst case, which can be improved to $O(gn^{2.5})$ under reasonable assumptions (nearest-neighbor chip connectivity, i.e., $d \in O(\sqrt{n})$, and an extended layer E with at most $O(n)$ CNOT gates).

It is important to note that the initialization step only needs to be repeated when the calibration data change but that requires to recover data from the Internet that can be a slow operation (on the order of several seconds).

B. INITIAL MAPPING

Heuristic-based mapping transition algorithms rely crucially on a good initial mapping to achieve the best results. A well-known algorithm when trying to approximate the global minimum of a scalar function with a discrete search space is simulated annealing. Simulated annealing is a metaheuristic designed to explore the search space by randomly selecting neighbors of the current state, evaluating them with the provided cost function and evolving in such a way that the algorithm will not be trapped into local minimums. The simulated annealing algorithm is depicted in Algorithm 2.

A modified version of simulated annealing has already been applied in [30], where a repetition parameter R is used to explore several neighbors at each temperature step. The authors consider a simple `get_neighbor` function that modifies randomly the current mapping π to a neighboring mapping π_{neighbor} . However, `get_neighbor` function is limited as it is not aware of the underlying hardware. This means that from the set of mappings generated by this function and evaluated by the simulated annealing procedure, several mappings can be excluded even before evaluating the mapping cost.

We aim to improve the initial mapping generated with the simulated annealing procedure by designing an HSA algorithm using an HA `get_neighbor` method to explore the neighboring mappings. To explore different mappings, we separate the `get_neighbor` procedure in three algorithms governed by a top execution policy. This top layer policy decides which one of the three algorithms the `get_neighbor` method should execute to obtain a new mapping. The policy, we used randomly chooses which algorithm to use from the value of a random number.

The first algorithm, called `shuffle`, does not change the physical qubits involved in the current mapping but changes how they are mapped to logical qubits. The most straightforward algorithm that can be used for this task is a random shuffle—we list the physical qubits involved in the mapping, randomly shuffle them, and obtain a new arbitrary mapping with the same physical qubits.

The second algorithm, `expand`, does not change the mapping between physical qubits and logical ones but replaces one of the physical qubits involved in the mapping by another physical qubit that is not part of the mapping. Instead of a hardware-unaware `expand`, we use an `expand` algorithm that tries to avoid separating the physical qubits in the current mapping into two disconnected groups. Moreover, the

Algorithm 2: Simulated Annealing.

input : Initial mapping π_0 , Cost function C ,
Neighbor computation function
`get_neighbor`, Initial temperature T_{init} ,
Final temperature T_f , Temperature evolution
constant Δ

output: Best initial mapping found π_{opt}

```

1 begin
2    $\pi \leftarrow \pi_0$ ;
3    $\pi_{\text{opt}} \leftarrow \pi_0$ ;
4    $T \leftarrow T_{\text{init}}$ ;
5    $\text{cost} \leftarrow C(\pi)$ ;
6    $\text{cost}_{\text{opt}} \leftarrow \text{cost}$ ;
7   while  $T \geq T_f$  do
8      $\pi_{\text{neighbor}} \leftarrow \text{get\_neighbor}(\pi)$ ;
9      $\text{cost}_{\text{neighbor}} \leftarrow C(\pi_{\text{neighbor}})$ ;
10    if  $\text{cost}_{\text{neighbor}} < \text{cost}_{\text{opt}}$  then
11       $\text{cost}_{\text{opt}} \leftarrow \text{cost}_{\text{neighbor}}$ ;
12       $\pi_{\text{opt}} \leftarrow \pi_{\text{neighbor}}$ ;
13    end
14    if  $\text{cost}_{\text{neighbor}} < \text{cost}$  then
15       $\text{cost} \leftarrow \text{cost}_{\text{neighbor}}$ ;
16       $\pi \leftarrow \pi_{\text{neighbor}}$ ;
17    else
18      if  $\text{rand}() < \exp\left(\frac{\text{cost} - \text{cost}_{\text{neighbor}}}{T}\right)$  then
19         $\text{cost} \leftarrow \text{cost}_{\text{neighbor}}$ ;
20         $\pi \leftarrow \pi_{\text{neighbor}}$ ;
21      end
22    end
23     $T \leftarrow T \times \Delta$ ;
24  end
25  return  $\pi_{\text{opt}}$ ;
26 end

```

algorithm encourages rearrangement of qubits based on the figure of merit chosen (i.e., final state fidelity, circuit depth, execution time). In this algorithm, we consider that strongly connected qubits have high fidelity. The HA implementation aims to identify the qubits with the least and most connections. Moreover, based on our tests with qubit measurement operations, we find there is a huge source of errors. To account for these errors, we add weights to qubit to determine the best and worst qubits in terms of their measured fidelity.

The third algorithm used in the `get_neighbor` algorithm is called `reset`. Its purpose is to give the possibility to the simulated annealing algorithm to escape local minimums. This algorithm is needed because the first two algorithms `shuffle` and `expand` will likely explore only the close neighborhood of the current mapping and may not be able to escape a local minimum. To avoid being stuck, the `reset` algorithm tries to find a potentially good new initial mapping from a randomly chosen qubit, without considering the previously explored mappings. The algorithm starts

with a random qubit and expands the mapping by iteratively weighting all the qubits and adding the best qubit to the new mapping.

C. METRICS

To evaluate our solution and compare it to other algorithms, we use some metrics that are described in the following paragraphs.

The first metric is the success rate of the mapped quantum circuit on a given hardware. We define the success rate of a quantum circuit as the fidelity of the quantum state obtained at the end of the execution of this quantum circuit on the hardware. We estimate this success rate by executing the quantum circuit a large number of times (8192), counting the number of executions that gave the expected answer and dividing this number by the total number of executions. The expected answer is obtained by executing the quantum circuit on a simulator.

The second metric chosen is the additional number of CNOT gates. This metric is tightly linked with the total number of SWAP gates inserted.

The third metric is the total execution time of the circuit. As the execution time of each CNOT gate can be extracted from [40], we can estimate the overall execution time for a given circuit. This metric is important for several reasons. First, it shows the ability of the mapping algorithm to schedule gates in parallel when possible and how good is the algorithm at doing this. Second, it allows us to have an idea of the importance of decoherence noise in the computed fidelity. For each qubit, the execution time is computed by adding the total execution time of gate operations acting on it. The longest qubit execution time is selected to represent the total execution time of the quantum circuit.

IV. EVALUATION AND COMPARISON OF THE PROPOSED HA ALGORITHM

A. METHODOLOGY

All the benchmarks used are collected from the previous works [28], [29], [44]. They include several functions taken from RevLib [45] as well as quantum algorithms from a variety of domains, including optimization, simulation, quantum arithmetic, etc. They are well known in the community and given as quantum circuits written in the OpenQASM language [42].

We chose two quantum chips, `ibmq_almaden` and `ibmq_valencia`, available from the IBM Quantum experience website and one quantum chip `ibmq_tokyo`, which is not accessible currently but widely used by state-of-the-art algorithms. `ibmq_almaden` is a 20-qubit quantum chip. Its topology and characteristics are summarized in Fig. 1 and Table 1. `ibmq_valencia` is a 5-qubit chip depicted in Fig. 2(b). `ibmq_tokyo` is a 20-qubit virtual chip depicted in Fig. 7. We execute benchmarks on `ibmq_valencia` and `ibmq_almaden` to check cases when the mapped quantum circuit needs all the available qubits or only a small number of them. Moreover,

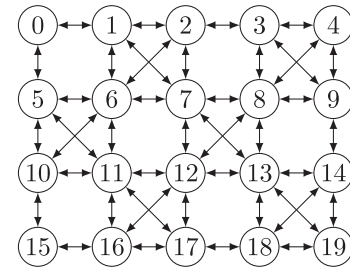


FIGURE 7. `ibmq_tokyo` topology.

we use `ibmq_almaden` and `ibmq_tokyo` to compare our algorithm with state-of-the-art algorithms in terms of number of additional gates. Note that, we do not have to execute the mapped quantum circuit on real quantum hardware to count the number of additional gates.

Our algorithm is implemented in Python and the Qiskit version is 0.19.1. To empirically evaluate our algorithm, we use a personal computer with one Intel i5-5300 U CPU and 8 GB memory. The Operating System is Ubuntu 18.04.

Several published qubit mapping algorithms are available as discussed in Section I. SABRE [29] seems to be the best algorithm at the time of writing when comparing the number of inserted gates to make the quantum circuit hardware-compliant. It provides a good initial mapping method and a mapping transition algorithm. Another algorithm dynamic lookahead (DL) [34] based on SABRE shows an improvement in terms of number of additional gates. Moreover, the mapping method presented in [19] uses the hardware calibration data to try to find a good mapping. We compare to all these algorithms. The source code of SABRE has been provided by the authors of the algorithm, and the mapping method presented in [19], called noise-adaptive (N-A) compiler, has been integrated into Qiskit as a transpiler pass. Finally, we also include the default transpiler included in Qiskit as the baseline. We execute our HA mapping transition algorithm with two different initial mapping algorithms—SABRE initial mapping algorithm and our HSA algorithm.

Summarizing, to test on real hardware, five different algorithms are included in the benchmarks: 1) our HA mapping algorithm with SABRE initial mapping; 2) our HA mapping algorithm with HSA initial mapping; 3) SABRE mapping algorithm with SABRE initial mapping; 4) N-A Compiler; and 5) Qiskit transpiler. For a fair comparison, we set the `optimization_level` parameter of the Qiskit transpiler to zero and make sure that the circuits obtained from the five methods are all executed with the same calibration data. The `optimization_level` is set to zero to invoke only the mapping transformation and not the optimization transformation. Moreover, when using the N-A compiler, the routing method is set to “lookahead” to make sure that it has the lookahead ability. To compare the number of additional gates without accessing to real hardware, three algorithms are included: 1) SABRE; 2) DL; and 3) HA.

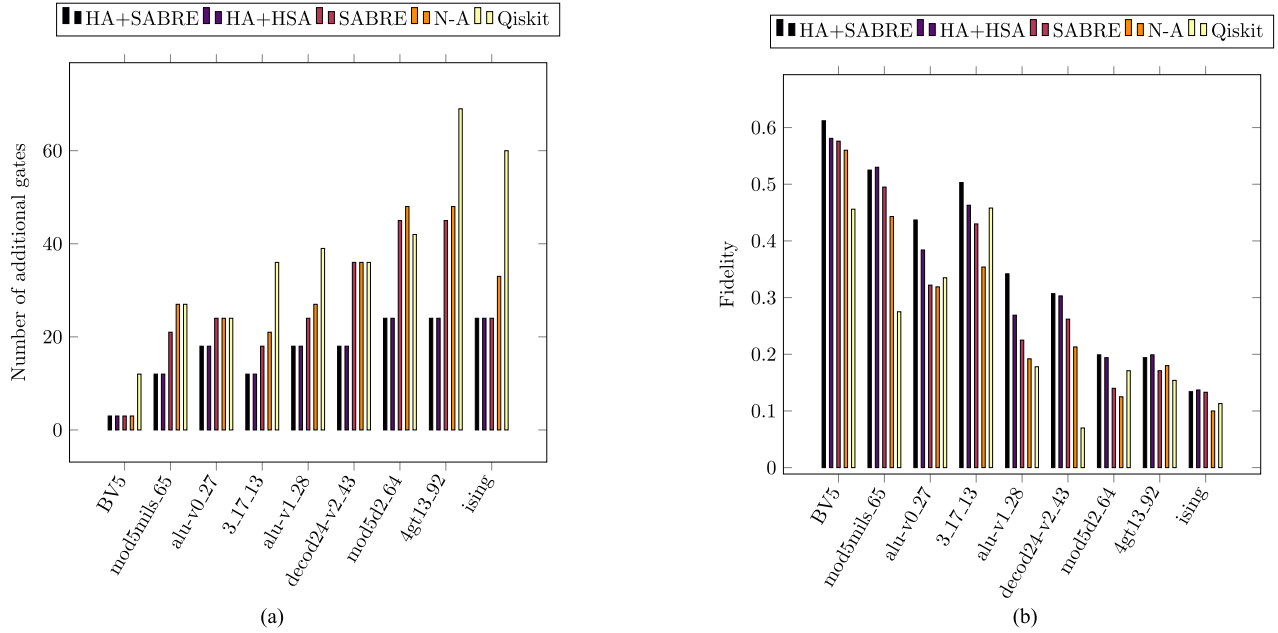


FIGURE 8. Comparison of number of additional gates and fidelity on `ibmq_valencia`. HA has been used with $\alpha_1 = 0.5$, $\alpha_2 = 0.5$ and $\alpha_3 = 0$. (a) Number of additional gates. (b) Fidelity.

To evaluate our algorithm with the different initial mapping methods, we allow each of them to call the mapping algorithm at most 100 times. The number of calls to the mapping algorithm is a natural parameter of the simulated annealing-based method, but the SABRE initial mapping method only needs two calls. To let the SABRE algorithm take advantage of a larger number of calls, we repeat the algorithm on several random initial mappings until no more calls are allowed and choose the best mapping found. The whole process is repeated 10 times to obtain 10 initial mappings.

We divide benchmarks by size according to their number of gates. We only execute small size benchmarks on real quantum hardware, because the other benchmarks with a large number of gate operations introduce too much noise to obtain any meaningful results. Moreover, the initial mapping generation process described above is applied on small and medium sized benchmarks. Large benchmarks suffer from long run time, so we generate 10 initial random mappings and use them with different algorithms. When using `ibmq_tokyo` virtual chip, we select the best results out of five attempts, which is a similar approach applied in SABRE and DL.

When testing the HSA algorithm, we used a random policy to choose which one of the three subroutines to execute. The `shuffle` procedure is executed with a probability of 0.9, the `expand` algorithm is chosen with a probability 0.08, and the `reset` procedure is executed when the two previous algorithms are not used (i.e., 0.02 probability).

First, we compare the number of additional gates and fidelity. The weight parameter α_1 of swap matrix S is set to 0.5, the weight parameter α_2 of CNOT error matrix \mathcal{E} is set to 0.5 and the weight parameter α_3 of CNOT execution time T is set to 0. Second, we compare the number of additional gates and

total execution time. Weight parameter α_1 of swap matrix S is set to 0.5, weight parameter α_2 of CNOT error matrix \mathcal{E} is set to 0, and weight parameter α_3 of CNOT execution time T is set to 0.5. Third, we compare the number of additional gates for circuits that are not executable on the real quantum device. Weight parameter α_1 of swap matrix S is set to 1, and the other two parameters are set to 0. For these three comparisons, the weight parameter W in the cost function is set to 0.5 and size of extended layer is set to 20.

B. EXPERIMENTAL RESULTS

We compare both the average number of additional gates [see Figs. 8(a) and 9(a)] and average output state fidelity [see Figs. 8(b) and 9(b)] among the 10 initial mappings for the five methods. The complete experimental results are listed in Tables 2 and 3.

The Qiskit default qubit mapping algorithm is nearly always the worst one in terms of additional gates, which translates in most of the cases to the worst output state fidelity. Although N-A Compiler takes into account the calibration data and has the lookahead ability, results show that it does not outperform the SABRE mapping algorithm with SABRE initial mapping (labeled as SABRE in the plots). Our HA mapping algorithm with SABRE initial mapping (labeled as HA+SABRE in the plots) seems to be the best combination as in average it achieves the best output state fidelity. Moreover, our HA algorithm with SABRE initial mapping gives the minimum number of additional gates. HA mapping algorithm with HSA initial mapping (labeled as HA+HSA in the plots) is also good, but its results are less consistent than HA+SABRE due to its random nature. Although, in many test cases, it outperforms SABRE.

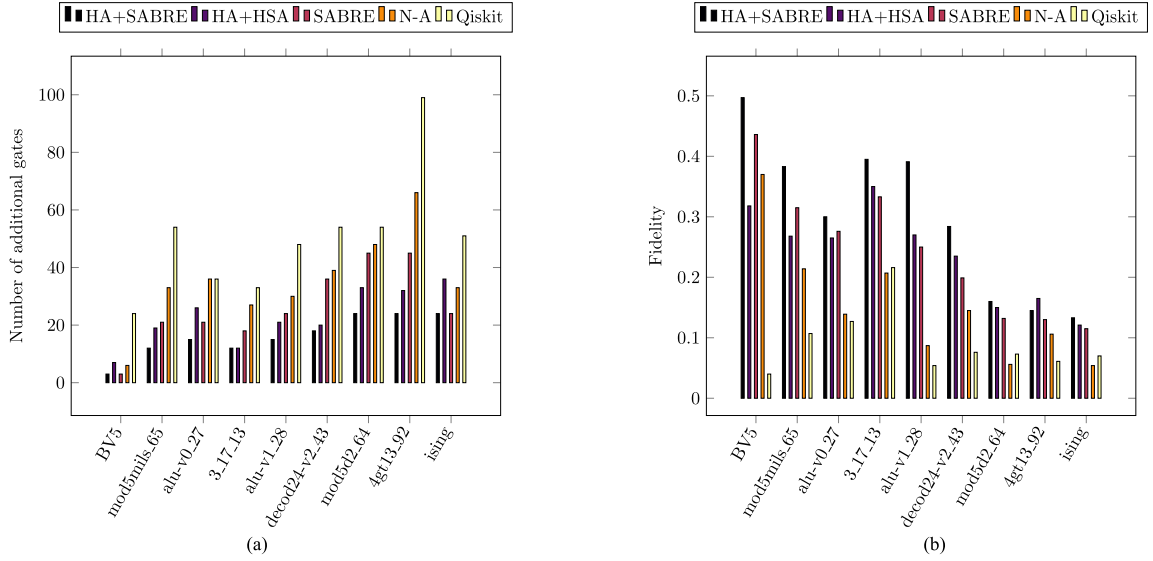


FIGURE 9. Comparison of number of additional gates and fidelity on `ibmq_almaden`. HA has been used with $\alpha_1 = 0.5$, $\alpha_2 = 0.5$, and $\alpha_3 = 0$. (a) Number of additional gates. (b) Fidelity.

TABLE 2. Comparison of number of additional gates and fidelity on `ibmq_valencia`

Original Circuit		SABRE					HA + SABRE					HA + HSA				Qiskit		N-A		Comparison			
name	n	g _{all}	g	g _{min}	S	S _{max}	g	g _{min}	S	S _{max}	t	g	g _{min}	S	S _{max}	g	S	g	S	$\Delta g\%$	$\Delta g_{min}\%$	$\Delta S\%$	$\Delta S_{max}\%$
BV5	5	15	3	3	0.576	0.639	3	3	0.612	0.639	0	3	3	0.581	0.63	12	0.456	3	0.56	0	0	6.3	0
mod5mils_65	5	35	21	21	0.495	0.515	12	12	0.525	0.559	0.003	12	12	0.53	0.559	27	0.275	27	0.443	42.9	42.9	6.1	8.5
alu-v0_27	5	36	24	24	0.322	0.329	18	18	0.437	0.437	0.002	18	18	0.384	0.431	24	0.335	24	0.319	25	25	35.7	32.8
3_17_13	3	36	18	18	0.43	0.476	12	12	0.503	0.546	0.004	12	12	0.463	0.542	36	0.458	21	0.354	33.3	33.3	17	14.7
alu-v1_28	5	37	24	24	0.225	0.233	18	18	0.342	0.384	0.004	18	18	0.269	0.384	39	0.178	27	0.192	25	25	52	64.8
decod24-v2_43	4	52	36	36	0.262	0.396	18	18	0.307	0.37	0.004	18	18	0.303	0.372	36	0.07	36	0.213	50	50	17.2	-6.6
mod5d2_64	5	53	45	45	0.14	0.208	24	24	0.199	0.207	0.005	24	24	0.194	0.207	42	0.171	48	0.125	46.7	46.7	42.1	-0.4
4gt13_92	5	66	45	45	0.171	0.191	24	24	0.194	0.206	0.006	24	24	0.199	0.22	69	0.154	48	0.18	46.7	46.7	13.5	7.9
ising	5	90	24	24	0.133	0.145	24	24	0.134	0.141	0.007	24	24	0.137	0.143	60	0.113	33	0.1	0	0	0.8	-2.8

HA has been used with $\alpha_1 = 0.5$, $\alpha_2 = 0.5$, and $\alpha_3 = 0$

n: number of qubits. **g_{all}**: total number of gates. **g**: average number of additional gates. **g_{min}**: minimum number of additional gates. **S**: mean of success rate. **S_{max}**: maximum of success rate. **Δg** : comparison of average number of additional gates between HA+SABRE and SABRE. **Δg_{min}** : comparison of minimum number of additional gates between HA+SABRE and SABRE. **ΔS** : comparison of mean of success rate between HA+SABRE and SABRE. **ΔS_{max}** : comparison of maximum of success rate between HA+SABRE and SABRE. **t**: runtime of HA+SABRE in seconds.

TABLE 3. Comparison of number of additional gates and fidelity on `ibmq_almaden`

Original Circuit		SABRE					HA + SABRE					HA + HSA				Qiskit		N-A		Comparison			
name	n	g _{all}	g	g _{min}	S	S _{max}	g	g _{min}	S	S _{max}	t	g	g _{min}	S	S _{max}	g	S	g	S	$\Delta g\%$	$\Delta g_{min}\%$	$\Delta S\%$	$\Delta S_{max}\%$
BV5	5	15	3	3	0.436	0.624	3	3	0.497	0.651	0.002	7	6	0.318	0.508	24	0.04	6	0.37	0	0	14	4.3
mod5mils_65	5	35	21	21	0.315	0.47	12	12	0.383	0.481	0.003	19	15	0.268	0.439	54	0.107	33	0.214	42.9	42.9	21.6	2.3
alu-v0_27	5	36	21	21	0.276	0.413	15	15	0.3	0.483	0.002	26	19	0.265	0.408	36	0.127	36	0.139	28.6	28.6	8.7	16.9
3_17_13	3	36	18	18	0.333	0.469	12	12	0.395	0.519	0.002	12	12	0.35	0.502	33	0.216	27	0.207	33.3	33.3	18.6	10.7
alu-v1_28	5	37	24	24	0.25	0.359	15	15	0.391	0.478	0.002	21	21	0.27	0.408	48	0.054	30	0.087	37.5	37.5	56.4	33.1
decod24-v2_43	4	52	36	36	0.199	0.334	18	18	0.284	0.401	0.006	20	18	0.235	0.387	54	0.076	39	0.145	50	50	42.7	20.1
mod5d2_64	5	53	45	45	0.132	0.198	24	24	0.16	0.266	0.003	33	33	0.15	0.263	54	0.073	48	0.056	46.7	46.7	21.2	34.3
4gt13_92	5	66	45	45	0.13	0.249	24	24	0.145	0.312	0.007	32	27	0.165	0.347	99	0.061	66	0.106	46.7	46.7	11.5	25.3
ising	5	90	24	24	0.115	0.177	24	24	0.133	0.191	0.01	36	30	0.121	0.235	51	0.07	33	0.054	0	0	15.7	7.9

HA has been used with $\alpha_1 = 0.5$, $\alpha_2 = 0.5$ and $\alpha_3 = 0$.

n: number of qubits. **g_{all}**: total number of gates. **g**: average number of additional gates. **g_{min}**: minimum number of additional gates. **S**: mean of success rate. **S_{max}**: maximum of success rate. **Δg** : comparison of average number of additional gates between HA+SABRE and SABRE. **Δg_{min}** : comparison of minimum number of additional gates between HA+SABRE and SABRE. **ΔS** : comparison of mean of success rate between HA+SABRE and SABRE. **ΔS_{max}** : comparison of maximum of success rate between HA+SABRE and SABRE. **t**: runtime of HA+SABRE in seconds.

We also tried to map and execute the `qft_10` circuit. We found that its output fidelity is less than 0.01 for all the methods tested in the benchmark. Because the base fidelity is too low to perform a meaningful comparison, we only compare the number of additional gates as summarized in Tables 4 and 5 for quantum circuits with a medium-to-large number of gates.

Fig. 10 shows the result of comparing the execution times, number of additional gates, and fidelities of our HA algorithm with SABRE algorithm on `ibmq_valencia`. The execution time is reduced by 19% on average. Even though the weight parameter α_2 of CNOT error matrix \mathcal{E} is set to 0, the fidelity is improved by 8%. The number of additional gates is reduced by 38%.

TABLE 4. Number of additional gates on `ibmq_almaden` for large circuits

Original Circuit				SABRE		HA			Comparison	
type	name	n	g_{all}	g	g_{min}	g	g_{min}	t	$\Delta g\%$	$\Delta g_{min}\%$
medium	qaoa	6	270	30	27	30	27	0.008	0	0
medium	ising_model_10	10	480	0	0	0	0	0.02	0	0
medium	ising_model_13	13	633	0	0	0	0	0.03	0	0
medium	ising_model_16	16	786	3	0	9	0	0.10	-200	0
medium	qft_10	10	200	93	81	66	42	0.04	29	48.1
medium	qft_13	13	403	192	177	195	171	0.07	-1.6	3.4
medium	qft_16	16	512	425	372	450	375	0.24	-5.9	-0.8
large	adr4_197	13	3439	2973	2856	2136	2004	2.13	28.2	29.8
large	radd_250	13	3213	2742	2655	2040	1926	1.62	25.6	27.5
large	z4_268	11	3073	2628	2559	1872	1815	1.44	28.8	29.1
large	sym6_145	14	3888	3024	2982	2022	1965	2.18	33.1	34.1
large	misex1_241	15	4813	3999	3831	2892	2630	3.04	27.7	31.3
large	rd73_252	10	5321	4539	4428	3261	3090	3.73	28.2	30.2
large	cycle10_2_110	12	6050	5127	5043	3795	3576	4.87	26	29.1
large	square_root_7	15	7630	6477	6324	4851	4707	7.00	25.1	25.6
large	sqn_258	10	10223	8679	8580	6012	5736	13.92	30.7	33.1
large	rd84_253	12	13658	11889	11673	8721	8574	24.54	26.6	26.5
large	co14_215	15	17936	16710	16368	13071	12426	37.81	21.8	24.1
large	sym9_193	10	34881	30558	30027	21900	21168	160.19	28.3	29.5
large	9symml_195	11	34881	30471	30129	21949	21168	151.84	28	29.7

HA has been used with $\alpha_1 = 1$, $\alpha_2 = 0$, and $\alpha_3 = 0$.

n : number of qubits. g_{all} : total number of gates. g : average number of additional gates. g_{min} : minimum number of additional gates. t : runtime in seconds. Δg : comparison of average number of additional gates between HA and SABRE. Δg_{min} : comparison of minimum number of additional gates between HA and SABRE.

TABLE 5. Number of additional gates on `ibmq_tokyo` for large circuits

Original Circuit				SABRE		DL		HA		Comparison
type	name	n	g_{all}	g	t	g	t	g	t	$\Delta g\%$
medium	ising_model_10	10	480	0	0.004	0	0	0	0.005	0
medium	ising_model_13	13	633	0	0.007	0	0	0	0.01	0
medium	ising_model_16	16	786	0	0.01	0	0	0	0.02	0
medium	qft_10	10	200	54	0.103	39	0.015	36	0.015	7.7
medium	qft_13	13	403	93	0.036	96	0.031	78	0.043	18.8
medium	qft_16	16	512	186	0.084	192	0.062	174	0.09	9.4
large	adr4_197	13	3439	1614	0.49	1224	0.218	882	1.41	27.9
large	radd_250	13	3213	1275	0.48	1047	0.186	840	1.24	19.8
large	z4_268	11	3073	1365	0.44	855	0.202	801	1.13	6.3
large	sym6_145	14	3888	1272	0.56	1017	0.202	786	1.71	22.7
large	misex1_241	15	4813	1251	0.89	1098	0.249	942	2.57	14.2
large	rd73_252	10	5321	2133	0.94	2193	0.343	1635	3.19	25.4
large	cycle10_2_110	12	6050	2622	1.35	1968	0.348	1719	4.02	12.7
large	square_root_7	15	7630	2598	1.5	1788	0.406	828	5.66	53.7
large	sqn_258	10	10223	4344	3.52	3057	0.563	2712	11.7	11.3
large	rd84_253	12	13658	6147	5.39	5697	0.892	3843	21.8	32.5
large	co14_215	15	17936	8982	9.51	5061	1.062	6429	36	-27
large	sym9_193	10	34881	16653	30.17	13746	2.091	11553	138.3	16

HA has been used with $\alpha_1 = 1$, $\alpha_2 = 0$, and $\alpha_3 = 0$.

n : number of qubits. g_{all} : total number of gates. g : minimum number of additional gates. t : runtime in seconds. Δg : comparison of minimum number of additional gates between HA and DL.

Table 4 lists the result of the number of additional gates on `ibmq_almaden`. Using the selection of SWAP and Bridge gate, our HA algorithm can outperform SABRE on circuits with different sizes. For medium circuits, HA gives similar results as SABRE and an improvement from SABRE for only one circuit among the eight circuits tested. For large circuits, HA outperforms SABRE and consistently reduces the number of additional gates by 28% on average. Table 5 shows the number of additional gates on `ibmq_tokyo` when comparing our HA algorithm with SABRE and DL.

DL outperforms SABRE and our HA algorithm can further reduce the number of additional gates by 14% on average. SABRE and DL only provide their runtime on `ibmq_tokyo`, the difference between runtime of the three algorithms is shown in Table 5. Note that, DL is written in C++ and tested on a normal personal computer. SABRE is written in Python and tested on a server with 2 Intel Xeon E5-2680 CPUs (48 logical cores) and 378 GB memory. Since there is an intrinsic speed difference between C++ and Python as well as the different devices used, the

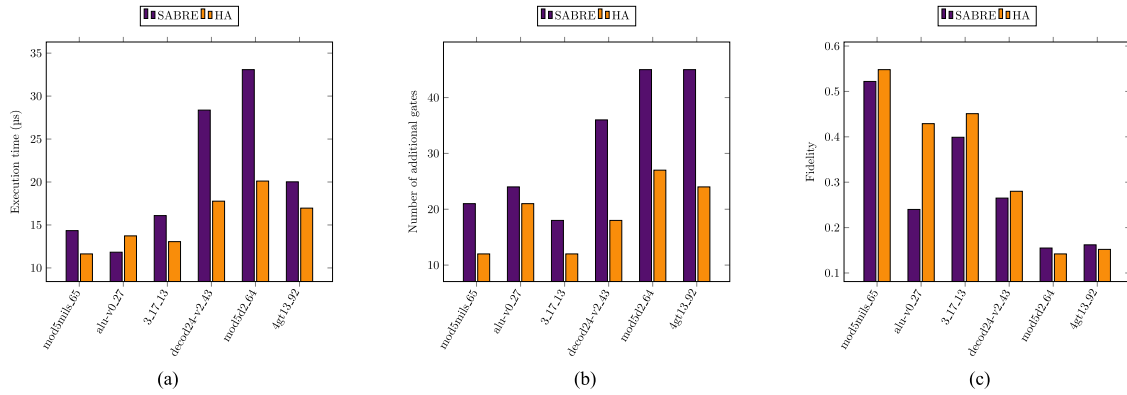


FIGURE 10. Comparison of execution time, number of additional gates and fidelity on ibmq_valencia. HA has been used with $\alpha_1 = 0.5$, $\alpha_2 = 0$, and $\alpha_3 = 0.5$. (a) Execution time. (b) Number of additional gates. (c) Fidelity.

runtime data in this table are for reference rather than for comparison.

V. DISCUSSION

A. DESIGN GUIDELINE

Given the current available NISQ hardware, it is important to adapt quantum programs to execute on such hardware while taking into account their physical constraints and limitations (noisy operations, number of qubits and gates). Here, we list several guidelines that can help a programmer to design quantum circuits that comply on given quantum hardware.

- 1) Check the topology and the calibration data of the device targeted. Try to map the most used qubit of the mapped circuit to the physical qubit that has the strongest coupling connection.
- 2) Try to apply a CNOT gate on qubits that are directly connected and with a reliable (i.e., low error rate) interconnect, so that no more additional gates are needed, and the overall circuit fidelity is improved.
- 3) If a CNOT gate cannot be applied on two-adjacent qubits, try to apply on two qubits, whose distance is two on the coupling graph. In such situation, one can select between a SWAP and Bridge gate to execute the CNOT gate. Also, the number of additional gates will be reduced.

B. FUTURE WORK

In this article, we present an efficient HA mapping algorithm based on heuristic search. For future studies, we find that the following potential research directions can be explored. First, our HA algorithm only takes into consideration the calibration data, which includes the gate error and the execution time. However, other physical constraints, such as crosstalk error may be included to take into account crosstalk coupling between interconnects. Second, we would like to investigate the adaptation of such a mapping algorithm to a multiprogramming mechanism as introduced in [46]. Executing multiple quantum circuits on the same chip allows us to use more efficiently hardware resources but may decrease the fidelity of the quantum operations due to unwanted interactions.

Finally, we find it relevant to investigate mapping algorithms for specific use cases such as quantum circuits constructed for quantum chemistry computations with VQE [13] or to solve linear systems with the VQLS algorithm [7].

VI. CONCLUSION

The quantum computers are now in the NISQ era. There's a gap between the design and execution of a quantum circuit in NISQ hardware. In this article, we present a HA heuristic for qubit mapping problem that adapts the quantum circuit to the quantum hardware. We design a mapping transition algorithm that uses calibration data and selects from either a SWAP or Bridge gate for qubit movement. Experimental results show that our algorithm can outperform state-of-the-art algorithms in terms of the number of additional gates, fidelity, and execution time. Our algorithm is evaluated on IBM quantum devices, but should be general enough to be used on quantum devices from other vendors as well.

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SUPPLEMENTARY INFORMATION

Authors have made available the source code, and it can be found at the following link: <https://github.com/peachnuts/HA>.

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