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Spatial Complexity of Reversibly Computable DAG

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ABSTRACT

In this paper we address the issue of making a program reversible in terms of spatial complexity. Spatial complexity is the amount of memory/register locations required for performing the computation in both forward and backward directions. Spatial complexity has important relationship with the intrinsic power consumption required at run time; this was our primary motivation. But it has also important relationship with the trade off between storing or recomputing reused intermediate values, also known as the rematerialization problem in the context of compiler register allocation, or the checkpointing issue in the general case. We present a lower bound of the spatial complexity of a DAG (directed acyclic graph) with reversible operations, as well as a heuristic aimed at finding the minimum number of registers required for a forward and backward execution of a DAG. We define energetic garbage as the additional number of registers needed for the reversible computation with respect to the original computation. We have run experiments that suggest that the garbage size is never more than 50% of the DAG size for DAGs with unary/binary operations.

Categories and Subject Descriptors

D.3.4 [Programming languages]: Processors—*Compilers*; F.1.2 [Computation by abstract devices]: Modes of computation—*Parallelism and concurrency*

General Terms

Design, Theory

Keywords

Reversible computing, Spatial complexity, garbage minimization

1. INTRODUCTION

One of the most important barriers to the performances of processors is power consumption and heat dissipation. Several works consider how to make architectures and programs

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cheaper in terms of energy. Thermal models are designed based on underlying electronics for architectures, or based on some software rules for programs, for instance balancing resources usage [23], or minimizing caches misses, or cutting unused devices. There is also a radically different approach that tackles energy issues under the point of view of intrinsic thermodynamics of computation and the basis argument that likewise thermodynamics transformations, irreversible programs have to dissipate heat. This originates from the Landauer remark [17] that erasing or throwing away a bit information must dissipate at least $kT \ln 2$ of energy. Therefore only reversible programs are likely to be thermodynamically adiabatic. Reversibility means here that no information is ever lost, it can always be retrieved from any point of the program.

Reversibility of programs has then been studied by Bennett [2]. A first easy way to make a program reversible is to record the history of intermediate variables along the execution, but then the issue of erasing - forgetting - that information, called garbage, remains. That garbage can be used as an estimation of the intrinsic energy consumption of programs and minimizing it is the objective. Bennett [2] proved that if the input can be computed from the output then there is a reversible way of computing the output from the input while eliminating the garbage. This may be at the price of large space usage for storing all the intermediate states during the computation. The alternative is to use checkpoints where the state of computation is stored and therefore from which the computation can be restarted for computing subsequent results. This is the very classical issue of trade-off between data storage and data recomputing in several domains of computer science.

It should be noted that evaluating energy with the size of garbage is not natural as one would better tend to relate energy with time complexity. Matherat et al. [20] elaborate on this point. They suggest first that convergences in state automata are the major sources of information loss and therefore heat dissipation. Second in the same idea as convergences they insist on the role of synchronisations in heat dissipation.

Reversibility of computation has several applications, among which we can mention bidirectional debuggers [8, 15], roll-back mechanisms for speculative executions in parallel and distributed systems, simulation and error detection techniques [6]. There are also algorithms that require a pass

where intermediate results have to be scanned in reverse order. This happens for instance in reservoir simulation, or for instance in automatic program differentiation [12]. Moreover reversible processor architectures have already been designed [14, 10], as well as reversible programming languages [10, 18].

In this paper we analyze the reversibility of computations with two aims. First we want to characterize the intrinsic reversibility of a program or piece of program based on its data dependency graph and not on some linear sequence of instructions. We hope this helps us understanding where energy bottlenecks arise from in ordinary programs. Second, because reversibility is very close to the trade-off between data storage and data recomputation, we want to analyze whether reversing operations or a set of operations may make the compiler problem of data rematerialization easier. An important hypothesis in this paper is that the operations that we consider can be reversed, this may concretely cause several problems that we skim over. We consider only DAGs (directed acyclic graphs), this means basic blocks in DDG (data dependency graphs). Compared to other works we consider possible rescheduling of instructions instead of a fixed sequence of already scheduled instructions. And the question that we address is: "Given a DAG computation graph with **reversible** operations, what is the minimum amount of garbage necessary to make the whole DAG reversible?"

Our ambition is therefore modest compared to this major and important issue of understanding precisely the relationship between physics, information, measurement, observers in one hand and information theory, computing in the other hand [11]. We believe however that this work may help optimizing scheduling and data storage in applications mentioned just before.

The remaining part of the paper is organized as follows. In Section 2 we explain more precisely our basic hypothesis and how we relate the problem of garbage minimization to the problem of register allocation, namely the number of registers required for reversibly executing the computation DAG. We give a heuristic algorithm for finding the number of additional registers ("garbage") required. In section 3 and 4, we propose a lower bound based on the decomposition of DAG into elementary paths and we compare the values lifetime between reversible and irreversible computing. Systematic experiments (section 5) with our heuristic algorithm suggests that garbage is never more than $n/2$ where n is the operation count in the DAG. We conclude (section 6) by a discussion on related works as well as our future work on storage and data recomputing for the minimization of memory access.

2. COST OF REVERSIBILITY AND ALGORITHM

In this section, we present our approach and algorithm for computing the spatial complexity of reversing a DAG. We consider a DAG of operations, typically the Data Dependency Graph of instructions within a basic block, see the part (a) of figure 1. Nodes of the graph are instructions denoted by the name of the variable carrying the result. This makes sense as two different nodes need to be treated as

two different variables. We consider only unary and binary operations and we make the important hypothesis that they can be made **reversible**.

2.1 Reversible Operations

A boolean function $f(x_1, x_2, \dots, x_n)$ with n input boolean variables and k output boolean variables is called reversible if it is bijective. This means that the number of outputs is equal to the number of inputs and each input pattern maps to a unique output pattern. Based on that we make the very rough abstract approximation that the operations in the DAG are reversible in the following sense: for unary operations, they are bijective so that the operand is uniquely determined by the result. For example, consider the increment function, defined from the set of integers Z to Z , that to each integer x associates the integer $y := x + 1$. The inverse function is $x := y - 1$, easily determined from the result uniquely. For binary operations, this means that only one additional value beside the result is needed for recovering both operands from this result and this additional value. This is typically the case of the basic arithmetic operations, like addition $c := a + b$, where (a, b) can be retrieved from (a, c) or (b, c) by a simple subtraction. Hence the '+' operation is considered as having two operands and two results.

This is only an abstraction and we are aware of a number of flaws underlying the concretization of this assumption. For instance the multiply '*' operation needs at least one additional resulting bit for determining which of both operands was 0 if the result is 0. There are also data precision issues especially with floating point operations and round-off problems but we neglect them and count only the number of data as a measure of memory/register space. We could also consider the semantics of operations and transform the operations in order to minimize the space for storing intermediate results. This is for instance done by Burckel et al. in [7] where they show that the computation transforming n inputs into n outputs can be (reversibly) performed by using a storage space not greater than n . Therefore in our abstracted model, when executing a binary operation we have the choice of memorizing the first or the second operand or both, provided that the reverse operation is possible based on the result and memorized operands.

What we want to evaluate is the maximum size of storage needed to execute the operations of the DAG reversibly in a forward and backward execution, or in other words we want to minimize the history required for performing the DAG reversibly. The additional storage space required compared to a simple forward execution is our criterion of "**energetic garbage**". With this criterion we want to characterize the intrinsic energetic cost of the DAG. Compared to the Bennett strategy for minimizing the storage space with checkpoints we have the degree of freedom to choose the schedule of operations. Finally our problem is *finding a schedule that minimizes the register requirement in a forward and backward execution. The difference between the bidirectional execution and the forward execution is called the energetic garbage.*

As an illustration, consider the code segment shown in figure 1(a) with its corresponding pseudo-assembly code and dependence graph in which each node corresponds to a state-

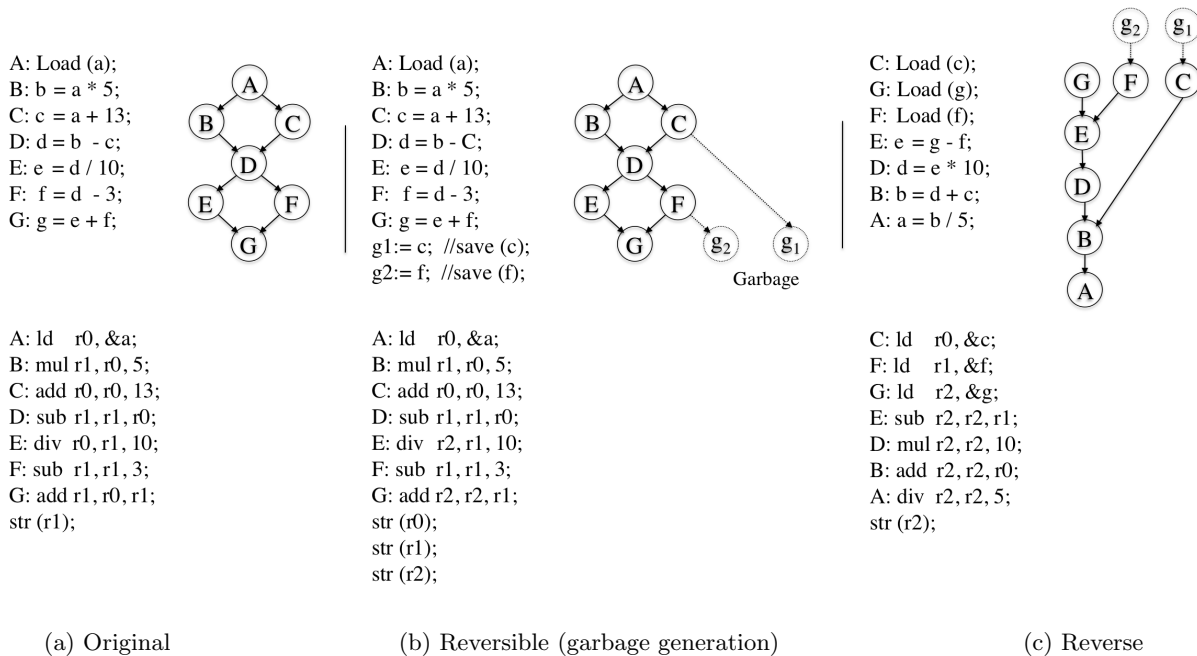


Figure 1: Illustrative example of reversible code and garbage generation.

ment in the code segment. This pseudo code requires only two registers. Figure 1(b) shows a reversible code that returns the same result as the code in the figure 1(a) and saves also two intermediate values. This code requires three registers for the three outputs g, f and c. Figure 1(c) shows the reverse code and its corresponding reverse graph derived from the reversible code in figure 1(b). It shows how a reverse computation could be performed to generate all previous values. Thus, 2 registers are required in the forward computation, 3 in the forward and backward computation. The energetic garbage is 1.

2.2 Algorithm

Like all optimization problems in register allocation it is likely that garbage minimization is a NP-complete problem. But we have not proven it. Here we describe a heuristic that schedules a DAG in the direct order while keeping some variables alive in order to make the backward computation feasible. We call garbage the difference between number of registers in the direct computation and number of registers in the schedule found in this algorithm.

Starting from the input data, the DAG is scheduled first in the direct direction and then in the reverse direction and we are looking for a schedule that uses the minimum number of registers. Since we consider a DAG, the number of registers required by a schedule is also the maximum number of simultaneously alive values during direct and reverse computation. One of the main issues of this work comes from the need to deal at the same time with the constraint of the minimum number of registers and with the constraint of saving values for enabling backward scheduling (reversibility constraint). Our algorithm decides two kind of informations: first which intermediate values will be saved in order to make the reverse computing feasible and second which successor node will kill the value - and hence will reuse the

same register in the actual register assignment of the direct computation.

The algorithm scans the DAG $G = (V, E)$ and schedules instructions in some topological order only in the direct direction. V is the set of nodes and E the set of edges. At each step the instruction with highest priority is selected. Algorithm is presented on the following chart.

Input: DAG computation graph.

Output: number of registers required for reversible computing

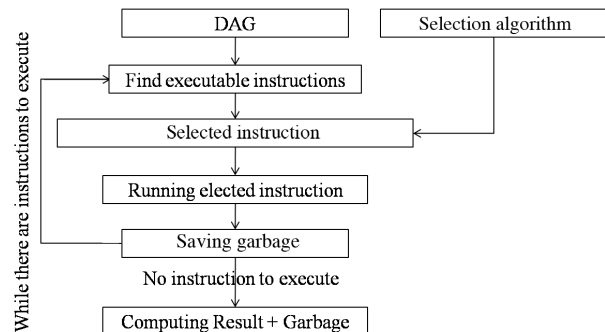


Figure 2: Scheduling algorithm diagram.

a. Priority

1. The order in which ready instructions are selected affects the number of registers required and the garbage size.
2. We use a heuristic.
3. Our scheduling is based on minimizing the number of resources (registers) without time constraint.
4. We favor instructions that have more predecessors with a minimum of successors in the DAG (this allows to use less registers)

5. We favor instructions on the critical path. This will increase the number of calls to the scheduler. We reuse only registers of the direct predecessors (to preserve information).

b. Labeling

Each node can take one of the following labels:

Active: if the value of a node is already computed and available in one of the registers.

Passive: any already calculated node that will not be used in a future computing.

Ready: if the value of a variable is ready to be calculated and all its predecessors are actives.

Idle: a node that is waiting because one of its predecessors is still waiting (waiting for all its predecessors to become active). Initially all source nodes (nodes without predecessors) are set to *Active*, the others are set to *Idle*.

Labeling rules

Let us consider the functions λ and Ω that define the state of a node u from the set of global nodes V

$$\lambda : V \longrightarrow \{Active, Passive, Ready, Idle\}$$

$$u \longrightarrow \lambda(u)$$

$$\Omega : V \longrightarrow \{Selected, Not\ selected\}$$

$$u \longrightarrow \Omega(u)$$

These rules are applied at each stage of calculation.

$\Gamma^+G(u)$ (resp. $\Gamma^-G(u)$) is the set of successor (resp. predecessor) nodes of node u .

$$\forall u \in V \wedge |\Gamma^-G(u)| = 0 \Rightarrow \lambda(u) = active$$

$$\forall v \in \Gamma^-G(u) \wedge \lambda(v) = active \Rightarrow \lambda(u) = ready$$

$$\exists u \lambda(u) = ready \wedge \Omega(u) = selected \Rightarrow \lambda(u) = active$$

Stop condition

$$\forall u \in V : \lambda(u) = active \vee \lambda(u) = passive$$

The final number of active nodes is the number of registers required.

The active nodes represent also the footprint needed for reversing the DAG computation.

Selection Algorithm

As an entry, the list of values ready to be calculated

Rule 1: u is the unique successor of v that remains to be scheduled: u will reuse the register used by v .

$$\text{if } \exists u \in V \lambda(u) = ready \left\{ \begin{array}{l} \text{if } \exists v \in \Gamma^-G(u) / |\Gamma^+G(v)| = 1 \left\{ \begin{array}{l} \Omega(u) = selected; \\ \lambda(v) = passive; \end{array} \right. \\ \left. \right\} \text{ else} \\ \text{if } \exists v \in \Gamma^-G(u) / \forall w \in \Gamma^+G(v) \wedge w \neq u \wedge \lambda(w) = active \left\{ \begin{array}{l} \Omega(u) = selected; \\ \lambda(v) = passive; \end{array} \right. \end{array} \right. \left. \right\}$$

Rule 2: Among ready nodes, one node with most successors is selected. For each pair of ready nodes $u, v \in V$

$$\exists u \forall v (u, v) \in V \lambda(u) = \lambda(v) = ready$$

$$\text{if } |\Gamma^+G(u)| \geq |\Gamma^+G(v)|$$

$$\Omega(u) = selected;$$

The second rule is applied if the first rule fails to select an instruction (node). Once we find the selected instruction, we go the labeling rules.

c. Guarantees

This algorithm ensures that all nodes will be covered, in

Nodes	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
Initialization	A	I	I	I	I	I	I	I	I	I	I	I	I	I	I	I
Labeling	A	R	R	I	I	I	I	I	I	I	I	I	I	I	I	I
Selection	-	-	S	-	-	-	-	-	-	-	-	-	-	-	-	-
.....																
Labeling	P	P	P	P	A	P	A	P	P	P	A	P	A	A	A	R
Selection	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	S
Labeling	P	P	P	P	A	P	A	P	P	P	A	P	A	P	A	A

A: Active P: Passive R: Ready I: Idle S: Selected

Table 1: An execution of the algorithm described by relabeling and selecting rules (corresponding to the code of example 2 figure 3)

other words, all values will be calculated. At the end active nodes contain active registers that will be used in the backward generation of all values.

d. Analysis

We can show that the following algorithm allows selecting a high priority instruction, which consumes less resources, at least at each computation step, this is a local optimization. At each computation step only one instruction is selected. The first rule allows the node who has a predecessor with one successor or predecessor of degree 1, to run first, this will not influence any other possible decision, and this node will reuse the register of this predecessor. The second rule is applied if the first rule fails to select an instruction. It consists of choosing the variable at the greatest distance from the result and which has more successors, the idea behind this is to increase the number of candidates from which the selected will be chosen. Therefore if we consider the number of candidates for the selection at step i is $NC_i = k$ then we want that at step $i+1$ $NC_{i+1} \geq k$. Labeling each calculated node as active will reduce the degree of successor nodes and increase the priority of neighboring nodes. An execution of this algorithm on the DAG of the figure 3(b) is shown in table 1. We will show that for every DAG G , the final configuration of registers allows the computation of all intermediate results back to input nodes.

e. Proof

We prove that from the nodes with active state we can find back all previous values.

At the end of the direct computation, we have only active or passive nodes, knowing that an active node changes its state to passive only during the call of the selection procedure and in the back track, an idle node becomes active iff: $\exists v \in \Gamma^+G(u) \lambda(v) = active \wedge \forall w \in \Gamma^-G(v) v \neq u \lambda(w) = active$

Proof by contradiction:

Assuming that there always exists a node with idle state $\forall instant t \exists u \in V \lambda(u) = idle \implies \forall v \in \Gamma^+G(u) \lambda(v) = idle \vee \exists v \in \Gamma^+G(u) \lambda(v) = active \wedge \exists w \in \Gamma^-G(v) v \neq u \lambda(w) = idle$

1. if $\forall v \in \Gamma^+G(u) \lambda(v) = idle$ by recursion we find that $\exists u \in V / |\Gamma^+G(u)| = 0 \lambda(u) = idle$ it's contradiction, because we know that : $\forall u \in V / |\Gamma^+G(u)| = 0 \implies \lambda(u) = Active$ (stop condition of the algorithm)

2. if one of its successors in active state which has another predecessor in the idle state: $\exists v \in \Gamma^+G(u) \lambda(v) = active \wedge \exists w \in \Gamma^-G(v) v \neq u \lambda(w) = idle$ this means there are two neighbor nodes in the idle state and each one

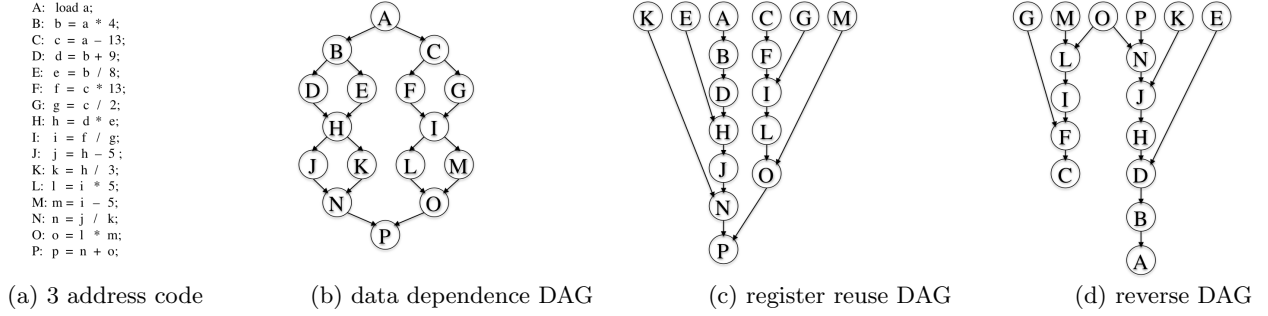


Figure 3: Example code and corresponding register reuse and reverse DAG

of them waits the other to become active, in other words both of them were passives at the end of the direct computation. A node becomes passive only during the call of the selection procedure; in this case we have either:

$|\Gamma^+G(u)| = 1 \wedge |\Gamma^+G(w)| = 1 \wedge \Gamma^+G(u) = \Gamma^+G(w) = \{v\} \wedge \lambda(u) = \lambda(w) = \text{passive}$. Which is impossible because v could not make two active nodes passive.

Let: $|\Gamma^+G(u)| = 1 \wedge |\Gamma^+G(w)| > 1$ This means that the state of w was modified from active to passive by another successor and not v , but as w is always waiting $\lambda(w) = \text{idle}$, it implies there is another idle neighbor that it can't modify its state and that there is another successor thanks to it, it took passive state previously.

Let take u_i the neighbor of u such that $|\Gamma^+G(u_i)| > 1 \wedge \lambda(u_i) = \text{passive} \implies \exists u_{i+1}$ neighbor of u_i and $|\Gamma^+G(u_{i+1})| > 1 \wedge \lambda(u_{i+1}) = \text{passive} \implies \exists u_{i+2} |\Gamma^+G(u_{i+2})| > 1 \wedge \lambda(u_{i+2}) = \text{passive} \implies \dots$ That leads us to the infinity, knowing that our graph is bounded (number of nodes is limited), which is a contradictory.

We have shown that our algorithm at the end of a direct computation allows to re-browse all of the graph in the reverse direction, that means recomputing all intermediate values. The table 1 shows the execution of the algorithm on the graph of figure 3.

3. REVERSIBLE DAG AND REGISTER REUSE DAG - LOWER BOUND

In this section we are seeking for a lower bound on the number of registers required for a DAG reversible computation. For that purpose we investigate the degree of register reuse in the reversible computation and we study its limitations in relation to the degree of dependency among variables.

Since the aim of reversible computing is the regeneration of data, some variables - contrary to the conventional computing - have to be saved even if they are not useful for the final result of (direct) computation. Therefore reuse is limited either between (a) independent values or (b) indirectly dependent values.

Register reuse between independent values. Independent variables correspond to variables for which no dependency path between both exist. They correspond to independent information. In contrast reuse chains [5] correspond to sequences of variables that can share the same register in some computation. Since we want to be able to

recover all intermediate values, this implies that for each reuse chain we need at least one live variable at each step of the computation. This means that the number of chains in a minimal decomposition is a lower bound for register requirement in reversible computing. This is actually also a consequence of the Dilworth theorem.

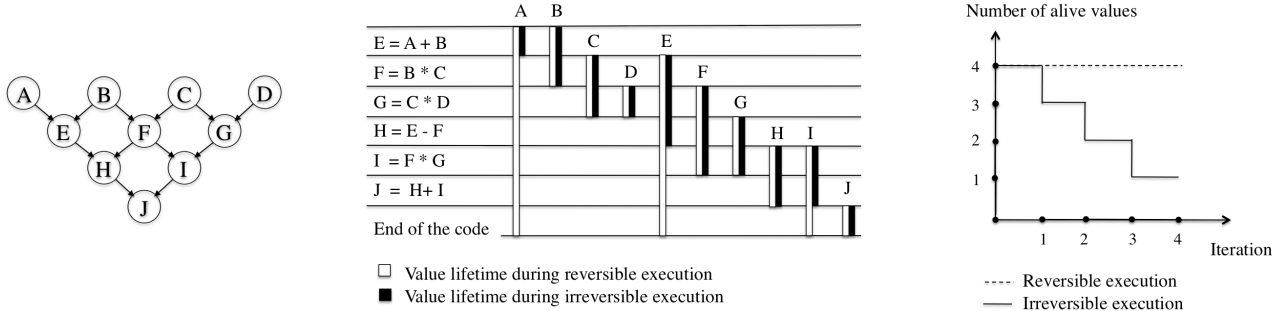
THEOREM 1. *The maximum number of independent elements in a partial order is equal to the number of chains in a minimal decomposition [9].*

For the graph of figure 3(b) this number is 4.

Register reuse between dependent values. For dependent values the main difference in reversible computation compared to direct ordinary computation is that we can not simply reuse the register of a killed value because of **convergences** in the graph. A convergence is a binary operation which has two operands and one result. This corresponds to loss of information that has to be stored for backward computation. This means for instance that stronger constraints for reuse have to be met for reuse chains than just ordinary chains. At least there must be an dependency edge between both operations in order that the second one can reuse the register of the first one. Therefore we need to consider decomposition of the graph into paths instead of chains. We prevent that a path contains a sub-path for the same reason (presence of a convergence).

The measurement of register requirement uses a Reuse DAG, which indicates which instructions can reuse a register used by a previous instruction. We use the same algorithm of the construction of the Reuse DAG for registers, proposed in [4], but the relation R that allows a value to reuse the register of a killed value requires that the defining instruction of the new value be the killing instruction of the previous one, in other words, it should be the last instruction that uses it. Therefore all allocation chains of the reuse DAG are paths in the original DAG.

In figure 3(a) we show the 3 address code for a basic block, the data dependence DAG using statement labels for the code in figure 3(b), and the register reuse DAG in figure 3(c), and reversible computing DAG in figure 3(d). In this example, at most four instructions can be executed in parallel because the number of chains in the minimal decomposition



(a) data dependence DAG

(b) values lifetime

(c) number of live values at each iteration

Figure 4: Example comparing number of live values and their lifetime between reversible and irreversible execution

is four. The values computed by independent statements D, E, F and G can all be alive at the same time and thus cannot share registers. Likewise, the values computed by J, K, L and M couldn't share registers since they are in separate chains, so there is no register reuse between K and M or K and G -contrary to the direct computation, nevertheless they are independent and they haven't the same live range. G and M are indirectly dependent, at the point of use of M, G is already dead and could normally share the same register. But we prevent this, because we need G and I for recomputing F. In this example there are five convergences in the graph, which means a loss in information at five points during computing, so we have to save additional information which is one of the inputs at each convergence, we can determine a lower bound of the size of this information by finding the minimal number of paths in the minimal decomposition of the reuse DAG. The sets of nodes $\{A, B, D, H, J, N, P\}$, $\{C, F, I, L, O\}$, $\{E\}$, $\{K\}$, $\{G\}$ and $\{M\}$ are all paths in the graph, and each end of each path can define an information that should be saved for the backward computation. The set of these paths is called register-reuse DAG because all nodes of a path are assigned to the same register. The reuse DAG chains those values that are not simultaneously live and can thus share a register, without any violation of information. Different paths are assigned to different registers, and therefore the number of paths is the number of registers. Thus, the register reuse DAG in figure 3(c) requires at least six registers. Therefore, by Theorem 2 a minimum decomposition of a DAG into elementary paths that don't contain any sub-path gives a lower bound to the register requirement for a reversible execution.

THEOREM 2. *The minimum number of registers required for a reversible execution of a program is bigger or equal than the number of elementary paths in a minimal decomposition of the corresponding DAG.*

4. REVERSIBILITY AND VALUES LIFETIME

There are cases when conventional and reversible computations consume the same number of registers, but values lifetime are not the same. As an example, consider the DAG in figure 4(a). Both of the executions -reversible and non-reversible, require 4 registers for the computation. A regis-

ter is used to hold a value from the time that the defining instruction executes until the value is killed by the last instruction that uses it. Figure 4(b) shows how a value that should be killed and frees a register, stays for the whole computation.

We try to compute the number of live values at every computation step. Since we allow only one instruction to be executed at each computation step, the curve is either constant or increasing in the reversible computing, figures 4(c) and 5; because we shouldn't get rid of a value if it's not directly replaced by another otherwise we will lose unrecoverable information.

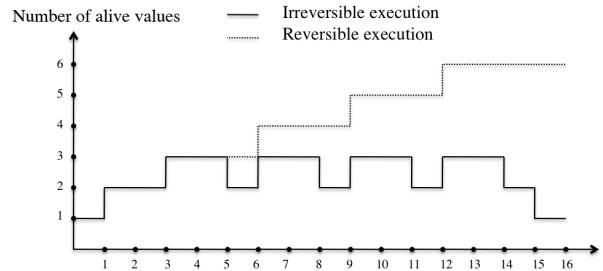


Figure 5: Example comparing the number of live values at each iteration step for the same code in reversible and irreversible execution (corresponding to the code of example 2 figure 3)

This can define another problem in the reversible computing: How to keep the minimum garbage with a shortest lifetime possible? We can deduce the minimum number of registers required for a reversible execution of a program from a direct execution by computing the number of live values at each step. The idea is to add the positive difference of the number of live values between two successive iterations in each computing step. If we take Ψ_i the number of live values at the iteration i with $i \in [1, N - \Psi_0]$ and N is the graph size or the number of nodes and Ψ_0 is the number of sources in the graph, q is an integer. This simple code can deduct the number of registers required for the reversible execution.

$$q = \Psi_0;$$

$$if(\Psi_{i+1} - \Psi_i > 0)$$

$$q+ = \Psi_{i+1} - \Psi_i;$$

In the reversible computing, we should not kill a datum if we have no other way to recompute it unlike in conventional computing. However, each increase of the number of live registers means that new values are generated and no one was killed. The possibility to clean garbage data (garbage-free) to free registers, can be done by a backtrack to recompute intermediate values or the input data.

5. EXPERIMENTAL RESULTS AND UPPER BOUND FOR THE GARBAGE SIZE

We come back to our algorithm defined in section 2. In order to understand more about what this criterion of garbage is, we made two kind of experiments. First we generated all graphs of some (small) fixed given size and second we used a randomly generated set of larger graphs. For exhaustive or random generation of DAGs we represent DAGs by their adjacency matrix, namely a boolean matrix with rows and columns labeled by graph vertices, with a 1 or 0 in position (v_i, v_j) according to whether v_i and v_j is an edge in the DAG or not. In both cases we computed the garbage as explained in previous sections.

From our experiments of small graphs we exhibited critical graphs for which garbage is maximal. These are the graphs of figure 6(b), where indeed according section 3 garbage is at least $n/2$ because a partition into elementary paths results in at least $n/2$ paths if the DAG has n nodes.

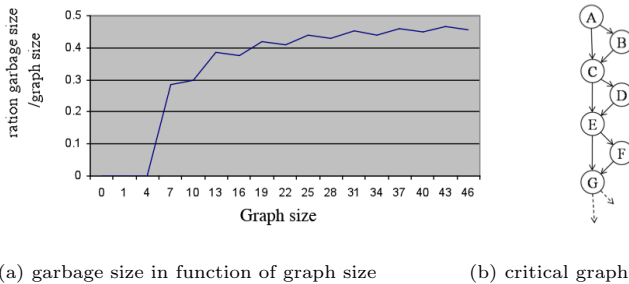


Figure 6: upper-bound to the garbage size

In [20] where reversibility of automata instead of DAG is considered, they argue that energetic garbage is related to convergences in automata - that result in loss of information. In this case our critical graphs do not have the maximal number of convergences. Maximal number of convergences correspond to DAG with only binary operations. In our case if we want to maximize the garbage size we have to maximize the number of convergences without increasing the number of register requirement in the direct computation. At least two registers are needed to create a convergence in a graph, so we fix the register requirement to two, and we try to create a maximum of convergences. We can observe that each increase of the number of convergence of one corresponds to an increase of the graph size of two, which explains the upper-bound of the garbage size (50% of the graph size).

This $n/2$ upper-bound is actually corroborated by experiments on randomly generated larger graphs. Figure 6(a) reports the maximum garbage obtained in percentage of the number of nodes. One can see that this number is never more than 50%.

Figure 7 reports also an histogram of garbage size for different size of graphs.

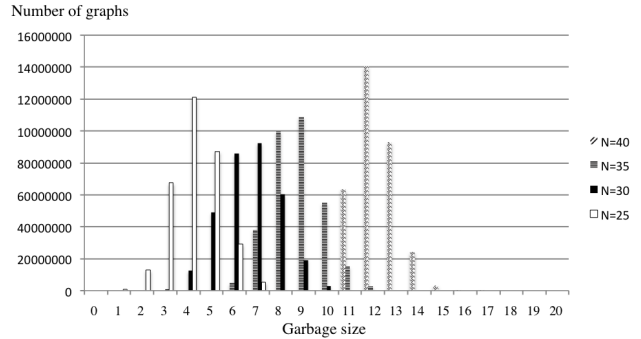


Figure 7: Number of graphs randomly generated according to the garbage size

6. RELATED WORKS AND CONCLUSION

Several works have addressed reversibility at the software level. There are basically two main approaches. Since writing reversible programs by hand is not quite natural, some works are devoted to the design of reversible programming languages [1], [18] Frank's R [10]. The alternative approach is to convert existing programs written in an irreversible programming language into equivalent reversible programs. An irreversible-to-reversible compiler receives an irreversible program as input and reversibly compiles it to a reversible program.

When one converts irreversible programs into reversible ones one has to face the issue of trading time complexity with data storage complexity. One elegant method was proposed by Bennett [3] where he models the former problem with a pebble game. Pebbles represent available data at some point of the program. One can add pebbles on some node when there is a way to compute that node with data identified by a pebble in a previous node. One can remove pebbles if there is an alternative way to recompute data required in this node. This is therefore an abstraction of reversible computations that allows analysis of the space and time complexity for various classes of problems, but this simulation operates only on sequential list of nodes. This sequence is broken hierarchically into sequences ending with checkpoints storing complete instantaneous descriptions of the simulated machine. After a later checkpoint is reached and saved, the simulating machine reversibly undoes its intermediate computation, reversibly erasing the intermediate history and reversibly canceling the previously saved checkpoint. Bennett chose the number of pebbles large enough ($n = O(\log T)$) so that m the number of steps become small. Other works such as [16] have also considered the reversible pebble game.

In [22], Vitanyi gives a time-space tradeoff for an irreversible computation using time T and space S to be simulated re-

versibly in time $T' = 3^k 2^{O(T/2^k)} S$ and space $S' = S(1 + O(k))$, where k is a parameter that can be chosen freely $0 \leq k \leq \log(T)$ in order to obtain the required tradeoff between reversible time T' and space S' . Naumann [21] considers the problem of restoring the intermediate values computed by such a program (the vertices in the DAG) in reverse order for a given upper bound on the available memory, while keeping the computational complexity to a minimum. He shows that the optimal data-flow reversal problem is NP-complete [21] but he doesn't consider reversible operations and seeks only reversible behavior for recovering intermediate values in reverse order. Griewank also in [12] presents an optimal time-space tradeoff algorithm in the context of automatic differentiation. In [19] it has been showed that the minimum number of garbage bits required to make a boolean function reversible is $\lceil \log(M) \rceil$, where M is the maximum of number of times an output pattern is repeated in the truth table. However, traditional techniques for bi-directional execution are not scalable to all classes of problems.

Conclusion

We have presented an analysis of the number of registers required to make a DAG computing graph reversible. We define the energetic garbage as the additional number of registers required for computing a forward and backward execution of the graph with respect to the simple forward execution. We gave a lower bound as the size of the decomposition of the graphs into elementary paths and through our experiments, we found that the garbage size does not exceed 50% of the program size - for DAG of unary/binary operations. However, values lifetime is shorter in a direct computation. A natural followup of this work would be to integrate this register allocation algorithm in a reversible compiler like the RCC -Georgia Tech's reversible C code compiler [13] in order to compare the size of generated codes with existing reversible programs and memory usage. But we also want to analyse more deeply in computation graphs the data rematerialization for recomputing data instead of reloading from memory - possibly after having spilled them in memory, in the presence and exploitation of reversible operations.

It is amazing to observe that in code optimization on current processors, minimizing power consumption amounts most of the time to minimizing the number of memory accesses, cache misses, etc. Therefore it was quite expectable that this thermodynamics view of computation leads to trade off between storage and recomputing. Understanding more thoroughly the relationship between variables of a computation graph in terms of mutual information between variables is our next step.

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