

Constructing the transition graphs of interacting hysterons

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Structure:

I. INTRODUCTION

Frustrated media can have rugged energy landscapes, with many metastable states. Recent work has shown that the many metastable states in frustrated systems can lead to unusual pathways when the system is subject to driving. For a variety of systems, including sheared amorphous solids[1][2], crumpled thin sheets[3], corrugated sheets[4] and origami structures[5] remarkable phenomena such as breaking of return-point memory, a multiperiodic response to cyclical driving, and transient memory as described by the park bench model [1][6] have been observed.

It turns out that the response of many frustrated systems can be captured by an interacting collection of hysteretic two-state elements, called hysterons. The hysteron model, first suggested by Ferenc Preisach[7], has long been used to model hysteretic systems as a collection of small hysteretic two-state elements. However, the Preisach model considers hysterons to be independent, while the interacting hysteron model has remained largely unexplored. The recent experimental findings show that the interacting hysteron model is relevant for many physical systems, and thus worth examining in more detail.

The possible metastable states and transitions between these states for a given hysteron system can be portrayed as a directed graph, which captures all possible responses of the system to driving[8]. Numerical sampling work from our group has shown that the number of possible transition graphs for a system of interacting hysterons increases dramatically with the number of hysterons. For as few as three interacting hysterons, over 15,000 distinct graphs are found[9]. Because the number of possible transition graphs quickly grows overwhelming, a systematic method will be needed for categorizing the graphs.

Here we establish a systematic method to gradually approach the complexity of the interacting hysteron model, starting from the Preisach model. We first establish a formalism for constructing the transition graph for coupled hysterons by starting from a Preisach graph, and adding the effects of coupling step-by-step. We show how this formalism is used to construct a list of candidate graphs for a given number of hysterons. We then detail a method for checking whether a given transition graph can be realized, both in general and for a specific model. Finally, we illustrate our formalism for a model where hysterons are linearly coupled.**REWORK ACCORDING TO CURRENT STRUCTURE**

II. MODEL

Here we detail the basic setup for simulating a system of N hysterons. We define the state of the system S as

$$S = \{s_1, s_2, \dots, s_N\} \quad (1)$$

where $s_i \in \{0, 1\}$ are the phases of individual hysterons. We then model the response of individual hysterons to an external driving field U by defining up and down switching fields U_i^+ and U_i^- . A hysteron transitions from phase 0 to 1 ('up') if $U \geq U_i^+$, and from phase 1 to 0 ('down') if $U \leq U_i^-$.

If the system of hysterons is non-interacting, then the switching fields $\{U_i^+\}, \{U_i^-\}$ are fixed values $u_i^{+,-}$. We refer to these switching fields in the absence of interactions as the bare switching fields. If there are interactions, the switching fields are modified by the state S :

$$U_i^{+,-}(S) = u_i^{+,-} + \Delta u_i(S) \quad (2)$$

When a driving field U is applied to a system with the collective state S , any hysterons that are in phase 0 and have $U \geq U_i^+$, or that are in phase 1 and have $U \leq U_i^-$, become unstable. If a single hysteron is unstable, that hysteron flips. If multiple hysterons are unstable, the system's response is ambiguous, as it is unclear which hysteron flips first and these operations do not commute if there is coupling. We therefore do not consider cases where multiple hysterons are unstable.

It can occur that the flipping of one hysteron destabilizes other hysterons when there is coupling. It may therefore take multiple iterations for the system to reach stability under a field U . We refer to this phenomenon as an avalanche. It is also possible that an avalanche leads back into the state itself, in which case the system becomes caught in a self-loop. We consider this scenario to be ill-defined, as the system's state is caught in an infinite loop.**TALK ABOUT ILL-DEFINED LAST**

The system's response to a field U is thus obtained by comparing the field to the switching fields $\{U_i^+\}, \{U_i^-\}$ of the hysterons individually, flipping hysterons that are unstable, and repeating this process until there are no more unstable hysterons. The response is ill-defined if multiple hysterons are unstable at the same time, or if the system becomes caught in an infinite self-loop.**REPETITIVE**

The possible responses of a collection of N hysterons to driving are captured by a state transition graph, which

consists of all stable states S and all transitions between states. For a system of N hysterons, there are 2^N possible states. In particular there are two saturated states, where either all $s_i = 0$ or all $s_i = 1$. Each state S has a single up and down transition, except for the two saturated states, which only have an up and a down transition respectively. Therefore, a transition graph for a collection of N hysterons consists of a maximum of 2^N states and $2(2^N - 1)$ transitions. It can occur that not all of the 2^N states are included in the transition graph, because it is possible for a state to not be reachable, or to only be reachable as an unstable intermediate state of an avalanche. (...) We now make two considerations that affect the number of distinct graphs, namely exchange symmetry and Garden-of-Eden states.

A. Uniqueness considerations

Fig. 2a: Exchange symmetry

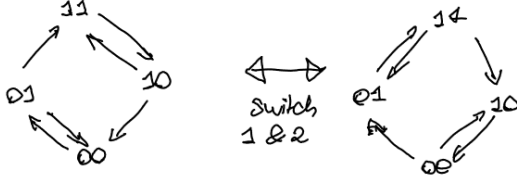


Fig. 2b: Garden-of-Eden states

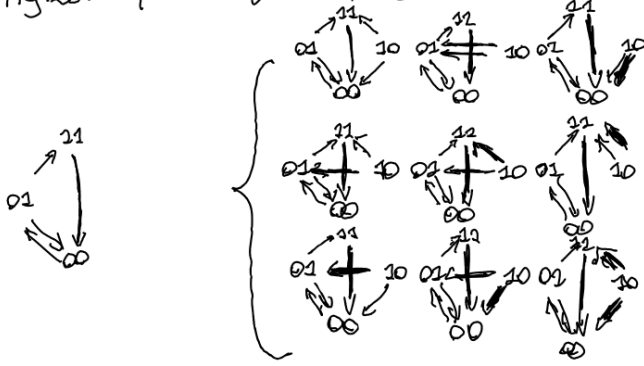


FIG. 1. Considerations for removing duplicate graphs: a) exchange symmetry and b) garden-of-Eden states.

We now describe two considerations that decrease the number of unique transition graphs, namely exchange symmetry and the existence of Garden-of-Eden states.

1. Exchange symmetry

When we define the state $S = \{s_1, s_2, \dots, s_N\}$, the labelling of hysterons 1 through N is arbitrary. As a

result, when two graphs only differ from one another by the ordering of the hysterons 1, 2, ..., N , we consider the graphs to be equivalent (Fig. 1a).

To account for exchange symmetry, we fix the order in which hysterons 1, 2, ..., N flip up when increasing the external field from the ground state. For example, for $N = 3$, we label the hysterons such that hysteron 3 flips up first from the state 000, hysteron 2 flips up first from 001, and trivially hysteron 1 flips up first from 011. Since there are $N!$ permutations of the order of N up switching fields, accounting for exchange symmetry decreases the number of candidate graphs by a factor $N!$.

2. Garden-of-Eden states

It often occurs that one or more of the possible states for N hysterons cannot be reached from the ground state. As is customary[5], we refer to states that cannot be reached as garden-of-Eden (GoE) states. We may choose to exclude GoE states in some cases, for instance in an experimental setup where one starts from the ground state. As a consequence of excluding GoE states, when two graphs only differ from one another by transitions that occur from GoE states, the graphs become duplicates. Therefore, when GoE states are ignored, the number of distinct graphs can turn out lower than the number of possible combinations of transitions.

B. Ordering of switching fields

In this section we discuss the crucial role of the ordering of the switching fields $U_i^{+,-}(S)$ for the topology of the corresponding transition graph. (...) **BRIEF OVERVIEW** Moreover, we discuss how the ordering of the switching fields provides an easy way to generate and label the transition graphs for non-interacting hysterons, while for interacting hysterons this is more challenging. (...) **CONCLUSION**

To establish the importance of the ordering of the switching fields, we consider how a transition graph is constructed for a given set of switching fields. **To start with, we ask which hysterons flip first from each state S , disregarding for the moment any avalanches that the initial hysteron flips might trigger?** **WHERE TO MENTION AVALANCHES?** To find the transitions from a state S , we vary the external driving U until it reaches a value where a hysteron flips, which we will call U^* . There are two possible transitions at each state, corresponding to an increase or decrease in the driving U [?]. For an increase in the driving U , the first hysteron to flip is that with the lowest up switching field, which we denote as

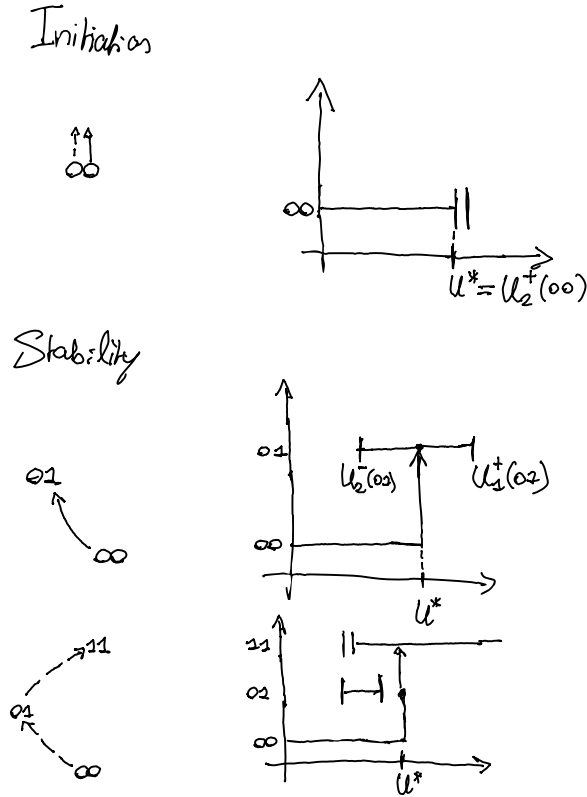


FIG. 2. Linear inequalities imposed by initiation & stability conditions of an example transition. **NO REFERENCE TO THIS FIG**

$U^+(S)$:

$$U^* = U^+(S) = \min_{s_j=0} U_j^+(S) \quad (3)$$

while for a decrease in the driving U , the first hysteron to flip is that with the highest down switching field, which we call $U^-(S)$:

$$U^* = U^-(S) = \max_{s_j=1} U_j^-(S) \quad (4)$$

We observe that the Equations 3 and 4 give conditions on the switching field which are linear inequalities, of the form $U_i^+(S) > U_j^+(S)$ and $U_i^-(S) > U_j^-(S)$. Notably, the linear inequalities of this form only impose an ordering between the switching fields $U_i^{+,-}(S)$. The values of the switching fields, as well as their relative distances, can be changed without changing the transition graph, so long as the ordering of the switching fields remains the same. Thus, to determine which hysterons can be flipped from each state, it is sufficient to know the ordering of the switching fields $\{U_i^{+,-}(S)\}$.

We now show that the ordering between the switching fields is sufficient not just to determine the first hysteron

flipped from each state, but to determine the entire transition graph, including any possible avalanches. To prove our statement, we look in detail at how a single transition is constructed. After the first hysteron has flipped, the system reaches a new state $S^{(1)}$. If the state $S^{(1)}$ is stable under U^* , then there is no avalanche. We call such a transition an elementary transition. If, however, the state $S^{(1)}$ is unstable, then an avalanche occurs and additional steps are required.

We first look at the case of an elementary transition, where the state $S^{(1)}$ is stable under the current driving U^* . Stability of $S^{(1)}$ under U^* means that all up switching fields $U_i^+(S^{(1)})$ are higher than the current driving U^* :

$$U_j^+(S^{(1)}) > U^* \quad (5)$$

while all down switching fields are lower than U^* :

$$U_j^-(S^{(1)}) < U^* \quad (6)$$

Unlike Equations 3 and 4, the stability conditions in Equations 5 and 6 do not make a direct comparison between switching fields. Rather, the switching fields at state $S^{(1)}$ are compared to the current driving, which is U^* . However, U^* itself is given by Equations 3 and 4, and therefore equal to one of the switching fields $U_i^{+,-}(S)$. Thus the stability of $S^{(1)}$ depends, once again, purely on the ordering of the switching fields.

Next we consider the additional steps that are required when $S^{(1)}$ is unstable. The system then enters into a new state $S^{(2)}$, for which we once again need to check the stability under the current driving. **ALSO OLD STABILITY CONDITIONS W/ ONE INEQ FLIPPED** The current driving is still held at U^* , so the stability conditions are as in Equations 5 and 6, except that the state $S^{(1)}$ is now replaced by $S^{(2)}$. Thus, if the state $S^{(2)}$ is stable, we obtain a set of linear inequalities as in Equations 5 and 6. If the state $S^{(2)}$ is unstable, then the system moves to a third state $S^{(3)}$, and in this way the system can keep iterating through states until a stable state is reached. **TOO QUICK** For each new state a corresponding set of inequalities is added, and each set of inequalities thus obtained only depends on the ordering of the switching fields. Hence, any transition $S \rightarrow S^{(1)} \rightarrow \dots \rightarrow S^{(l)}$, of any length l , is fully determined from the ordering of the switching fields. **INITIAL STATE S OR S⁽⁰⁾?**

Finally, we state that the full transition graph is described by a set of inequalities that is simply a combination of the conditions on the individual transitions. (...) We conclude that the corresponding transition graph for a set of switching fields only depends on the ordering of the switching fields, as described by a set of linear inequalities. We refer to the set of inequalities underlying a graph as the design inequalities[9].

In the case where hysterons are non-interacting, there are only $2N$ switching fields $\{u_i^{+,-}\}$. In addition, because we never compare an up switching field with a

down switching fields in the absence of avalanches, the only orderings that can change the graph are the partial orderings of the up switching fields $\{u_i^+\}$ and of the down switching fields $\{u_i^-\}$. The number of orderings that can lead to a unique graph is then $N! \times N!$. When we fix the up switching fields to account for exchange symmetry, we find that the number of permutations is only $N!$. Indeed there are $N!$ Preisach graphs for a collection of N hysterons, each corresponding to a different ordering of the down switching fields[10].

The ordering for each Preisach graph can be easily identified by considering the main loop: that is, the sequence of down transitions connecting the state with all $s_i = 1$ to that with all $s_i = 0$. We accordingly use the main loop as an identifier for the Preisach graphs. For instance, for $N = 2$ there are two Preisach graphs, corresponding to the main loops $(1, 2)$ and $(2, 1)$ [10].

We see that for the Preisach model, there is a one-to-one relationship between the permutations in the orderings of the switching fields $u_i^{+,-}$ and the possible transition graphs. When hysterons are allowed to interact, however, the situation changes. On one hand, there are then $2(2^N - 1)$ states $U_i^{+,-}(S)$, and because of the possibility of avalanches, the ordering between all of the states must be considered. On the other hand, the state transition graph is only determined by a partial order between some of the switching fields. As a result, the number of permutations of the switching fields $(2(2^N - 1))!$ for interacting hysterons greatly outnumbers the number of unique state transition graphs. For $N = 2$, the number of permutations is $8! = 40320$ while we know the number of graphs to be only 13. For $N = 3$, the number of permutations is as high as $24! \approx 6 \times 10^{23}$. The gross overestimation of the number of graphs makes it a challenge to find and identify all graphs for an interacting system of hysterons.

Even so, we can make use of the fact that the state transition graph only depends on a partial ordering of the switching fields. Namely, we know that each graph has an underlying set of linear inequalities $U_i(S_A) > U_j(S_B)$, which we refer to as the design inequalities[9]. The solution space of the design inequalities provides the region in parameter space that leads to the given graph. In particular, if the design inequalities have no solution, then the given graph is not realizable. **REWORK; SEE OUTLINE**

III. CONSTRUCTING POSSIBLE GRAPHS

In recent work, the transition graphs for two- and three-hysteron systems were found using numerical sampling. The obtained graphs contain several noteworthy phenomena that are not possible in the Preisach model, such as breaking of loop return point memory, scrambling, multiperiodic cycles and long transients[9]. A significant disadvantage of numerical sampling is

that it is not conclusive, as it is always possible that a region in parameter space leading to a certain transition graph is missed.**MOVE EARLIER** Here we aim to establish a tool that can conclusively find all possible graphs for a given number of hysterons N . We base our method on the design inequalities, as introduced in Section II B and elaborated on in the Supplemental Information. We use the design inequalities to argue that the transitions that make up a graph, whether they are single flips or avalanches, can only be constructed out of a fixed collection of elementary steps. We use our method to construct a list of candidate graphs, then filter out the realizable graphs by explicitly checking the solution space of the design inequalities.

A. Candidate graphs

A well-defined transition graph for N hysterons consists out of 2^N states, and $2(2^N - 1)$ transitions, namely the up and down transitions from each state, except for the saturated states which only have one transition each. A transition graph is constructed by choosing a transition path for each of these $2(2^N - 1)$ transitions. A transition path may be a single hysteron flip, or may consist of multiple flipped hysterons. When a transition path consists of multiple hysteron flips, we speak of an avalanche.**COMBINE W/ PREVIOUS SECTION** Naively, one can construct all possible transition graphs by finding the possible transition paths for each individual transition. For example, for $N = 2$ one finds $3^4 \times 6^2 = 2916$ graphs (Fig.3)**START W/ BASE GRAPHS**.

B. Base graphs

We now argue, based on the design inequalities, how avalanches must be constructed from a fixed choice of elementary steps.

For the hysteron i to flip up from the state S , the up switching field of hysteron i must be the lowest of all the hysterons:

$$U_i^+(S) = \min_{s_j=0} U_j^+(S) \quad (7)$$

Now suppose that the state S is not stable, but is instead a part of an avalanche. Let U^* be the field at which the avalanche is initiated. Hysteron i must then have

$$U_i^+(S) \leq U^* \quad (8)$$

while all other hysterons j with the state $s_j = 0$ must be stable, and therefore have

$$U_j^+(S) > U^* \quad (9)$$

Combining equations 8 and 9, we find that Equation 7 is implied.

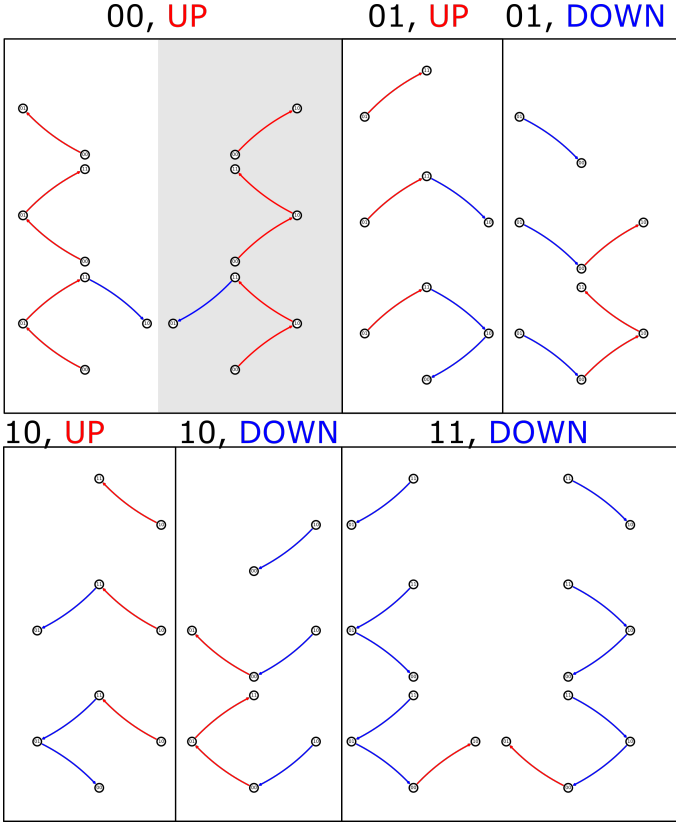


FIG. 3. Overview of the possible transition paths for each $N = 2$ transition individually. The greyed-out area signifies transitions that are left out when accounting for exchange symmetry.

(...)

C. Realizability

To better understand why some graphs are not realizable, we look in detail at the design inequalities underlying the graphs (see SI), and identify the inequalities that are in conflict with one another. By then tracing these inequalities back to the transitions from which we arise, we find the combinations of transitions (i.e., subgraphs) that are incompatible and thus make a graph not realizable. Several examples of impossible subgraphs are given in Fig. 4a. WHERE TO TALK ABOUT INCOMPATIBLE SUBGRAPHS?

We now direct our attention to a particular group of impossible subgraphs, an example of which is shown in Fig. 4b. The example subgraph consists of the single-flip transition $11 \downarrow 10$, and the avalanche $10 \uparrow 11 \downarrow 01$. We observe that the transition $11 \downarrow 10$ requires that $U_2^-(11) > U_1^-(11)$ to have hysteron 2 flip down from the state 11. Conversely, the transition $10 \uparrow 11 \downarrow 01$ requires the inequality $U_1^-(11) > U_2^-(11)$ to have an avalanche where only hysteron 1 flips down from the state 11. The

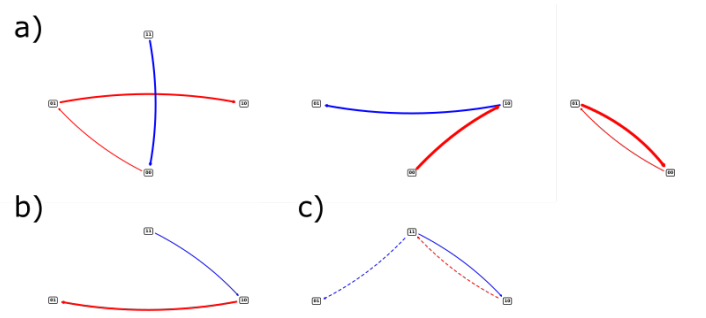


FIG. 4. Examples of impossible combinations of transitions for the two-hysteron system. a) Several examples of impossible combinations of transitions. b) Example of a combination of transitions that is impossible due to a violation of the underlying base graph. c) Illustration of how the combination of transitions in (b) violates the underlying base graph.

contradiction between the two transitions arises because the avalanche $10 \uparrow 11 \downarrow 01$ imposes that the switching field U_1^- is the highest down switching field, even though the transition $11 \downarrow 01$ is only a part of a larger avalanche. The argument used for the example subgraph can be extended to the general argument that the up and down flips which can occur from a state do not depend on whether or not that state is stable. As a consequence, all avalanches occurring in a graph must conform to an underlying graph of single-flip transitions, which we refer to as the base graph. PART OF BASE GRAPH ARGUMENT

Using base graphs, we can improve upon our method for generating candidate graphs. Namely, by first constructing all base graphs, and only then evaluating all possible avalanches for each base graph individually, we exclude the subgraphs of the type shown in Fig. 4b from the outset. Because the number of graphs for which the design inequalities need to be explicitly generated is then reduced, the base graph method has an improved efficiency compared to the naive approach.

In addition to the base graph method being an efficient method for generating candidate graphs, it also provides us with a powerful framework for understanding the effects of hysteron interactions on the allowed transition graphs.

D. The base graph method

1. Base graphs

We here show how we generate the list of base graphs for N hysterons, which is subsequently used to construct the full list of candidate graphs. We first explain how base graphs can be divided up into Preisach and scram-

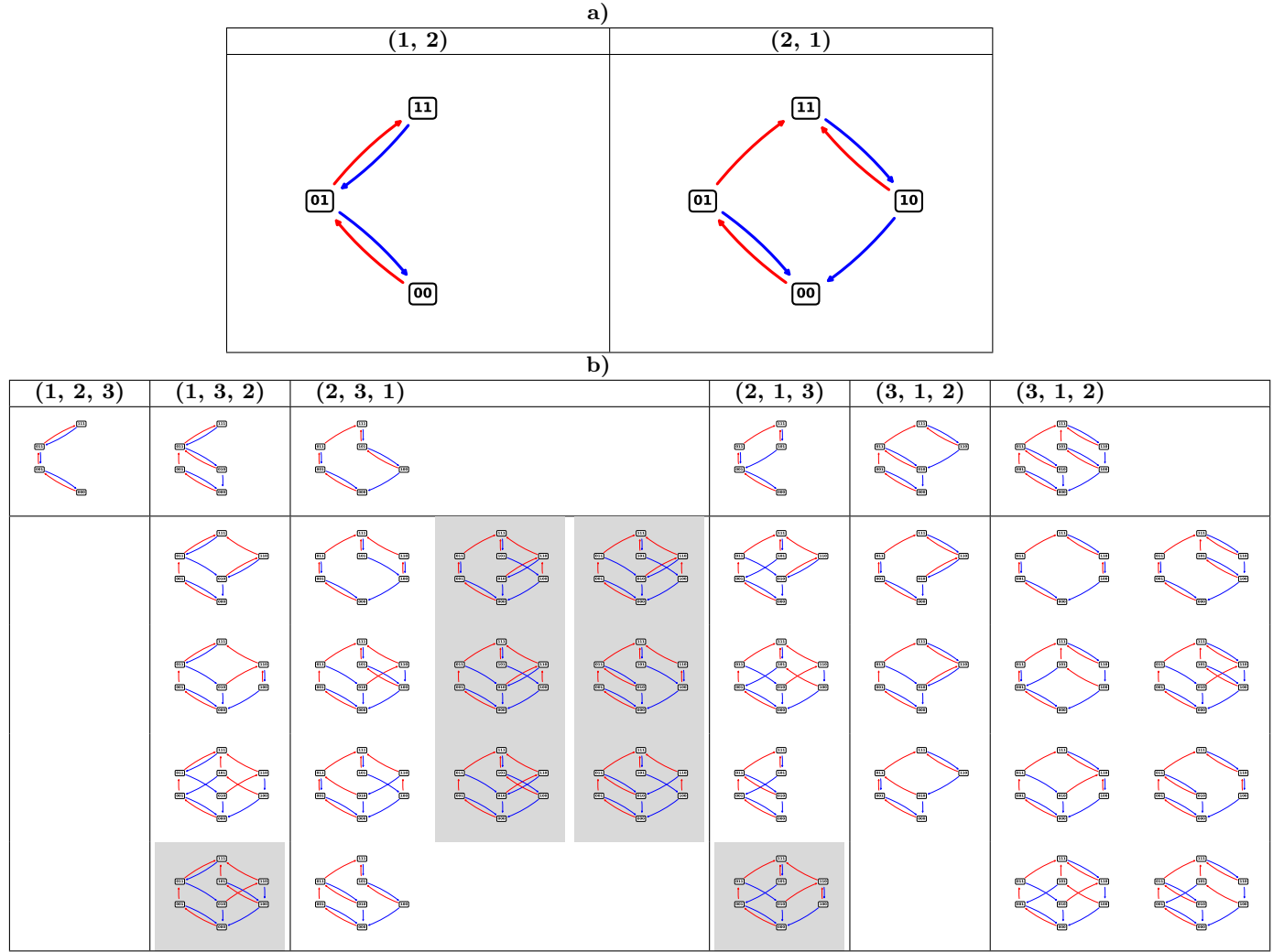


TABLE I. All possible base graphs for the two- and three-hysteron systems, indexed per main loop. a) All base graphs for $N = 2$; since scrambling is not possible for $N = 2$, the base graphs are the two Preisach graphs. b) All Preisach (top) and scrambled (bottom) base graphs for $N = 3$. Graphs on a gray background are possible for general, but not linear coupling.

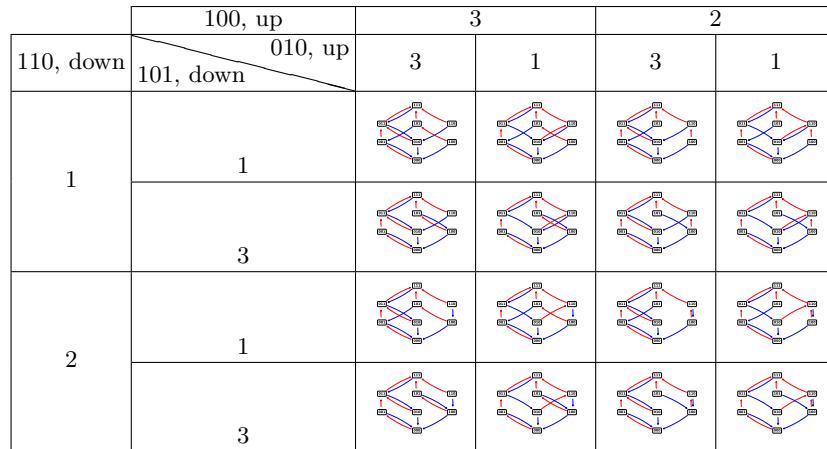


TABLE II. All sixteen base graphs for the example main loop (1, 3, 2). GoE pruning reduces the number of distinct base graphs from sixteen to four, out of which three are solvable in the linear model. **NO REFERENCE TO THIS FIG**

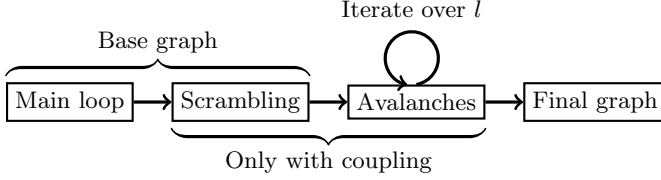


FIG. 5. Schematic of the construction of a graph via the base graph method. The base graph can itself be separated into the main loop and a number of scrambled transitions. **NO REFERENCE TO THIS FIG**

bled graphs. We then detail how one can count the number of base graphs per main loop.

We can distinguish between two groups of base graphs. First, there are the Preisach graphs, which are possible without coupling. **Second**, there is a group of scrambled graphs, which are possible for interacting hysterons. Scrambling is a phenomenon previously seen in interacting hysterons[9][4], where the order in which hysterons flip is not the same throughout the graph. It occurs when the ordering between the switching fields of two hysterons in the same phase are changed by one or more hysterons, and is therefore only possible for three or more hysterons. Remarkably, having introduced the concept of base graphs, we can identify scrambling as a feature of the base graph. We can thus separate the two coupling effects of scrambling and avalanches. **TOO QUICK; BUILD UP TO SCRAMBLING GRADUALLY**

In order to discuss scrambling more explicitly, we here define the main loop as the series of up and down transitions connecting the two saturated states. In the Preisach model, the ordering of the main loop fixes the transitions throughout the transition graph, so that each main loop corresponds to a unique Preisach graph. We additionally define the up and down boundary of the main loop as the series of up (down) transitions going from the state with all $s_i = 0$ (1) to that with all $s_i = 1$ (0). The number of main loops is then the number of permutations of the up boundary times that of the down boundary. Because both the up and down boundary consist of N flips, there are $N!^2$ main loops. However, we fix the up boundary when accounting for exchange symmetry, thus dividing the number of main loops by $N!$. Therefore, when accounting for exchange symmetry, there are only $N!$ main loops and resulting Preisach graphs. **REFER BACK TO ORDERING SECTION**

In a scrambled graph, one or more of the transitions in the graph deviates from the main loop, so that each main loop corresponds to a number of base graphs (Table.I). We now explicitly calculate the number of base graphs per main loop. To start with, we observe that the number of possible up and down flips for a state S depends on the number of hysterons that is in phase 0 or 1. Defining the magnetisation of a state as

$$M(S) = \sum_i s_i \quad (10)$$

we can state that the number of possible up flips for a state S is equal to $N - M(S)$, while the number of possible down flips is $M(S)$.

Next, we ask how many states exist with the magnetisation M . The number of states with the magnetisation M is the number of possible binary numbers that has M ones and $N - M$ zeros, which is equal to the binomial coefficient $\binom{N}{M}$.

It follows that the number of base graphs for a given number of hysterons N is

$$\#graphs = \left(\prod_{M=0}^N M \binom{N}{M} \right)^2 \quad (11)$$

When the up boundary of the main loop is fixed to account for exchange symmetry, the number of base graphs is reduced by a factor $N!$. When the down boundary is also fixed so that the main loop is fixed, the number of base graphs is reduced by another factor $N!$. Thus, one obtains the number of base graphs per main loop:

$$\#graphs/ML = \frac{\left(\prod_{M=0}^N M \binom{N}{M} \right)^2}{N!^2} = \left(\prod_{M=1}^{N-1} M \binom{N}{M}^{-1} \right)^2 \quad (12)$$

The number of graphs given by Equation 12 takes into account exchange symmetry, but not garden-of-Eden states (Section III A). When garden-of-Eden states are excluded, the number of base graphs turns out lower than that given in equation 12. For example, for $N = 3$, the number of base graphs provided by Equation 12 is 96, while excluding garden-of-Eden states reduces the number of base graphs to 35 (Fig. I).

2. Avalanches

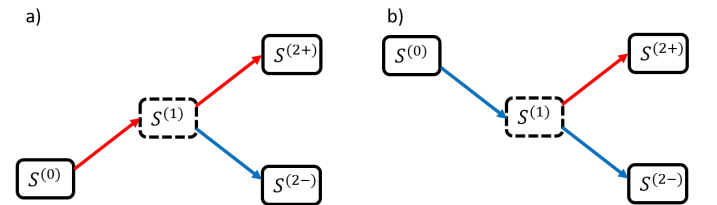


FIG. 6. Construction of an $l = 2$ avalanche by combining an incoming transition with one of at most two outgoing transitions. a) Combination of an incoming up transition with an outgoing up transition (ferromagnetic) or down transition (antiferromagnetic). b) Combination of an incoming down transition with an outgoing up transition (antiferromagnetic) or down transition (ferromagnetic)

We here show how the possible avalanches are evaluated for a given base graph. We first describe how one systematically constructs the possible avalanches of length $l = 2$ for a base graph. We then generalize our method

	00, UP	00→01		00→01	
11, DOWN	01, UP 10, DOWN	01→11	01→11→10	01→11	01→11→10
11→10	10→00				
	10→00→01				
11→10→00	10→00				
	10→00→01				

FIG. 7. All combinations of $l = 2$ avalanches for one of the two base graphs for $N = 2$ (top-left) graph, with indication of graphs that of which the design inequalities are solvable (green), solvable except for the final set of relaxation inequalities (orange) and unsolvable (red).

to find graphs containing avalanches of any length l .

To help our notation, we define L as the length of the longest avalanche in a graph. For example, $L = 2$ denotes a graph that contains at least one avalanche of length $l = 2$, and no $l = 3$ avalanches. In particular, $L = 1$ denotes a base graph.

Let a base graph contain the up transition $S^{(0)} \uparrow S^{(1)}$. From the state $S^{(1)}$ there are two possible transitions: an up transition $S^{(1)} \uparrow S^{(2+)}$, and a down transition $S^{(1)} \downarrow S^{(2-)}$. The transition $S^{(0)} \uparrow S^{(1)}$ can thus be extended to a $l = 2$ avalanche in two ways: either by the ferromagnetic avalanche $S^{(0)} \uparrow S^{(1)} \uparrow S^{(2+)}$, or by the antiferromagnetic avalanche $S^{(0)} \uparrow S^{(1)} \downarrow S^{(2-)}$ (Fig. 6a). Similarly, a down transition $S^{(0)} \downarrow S^{(1)}$ can be extended to the antiferromagnetic avalanche $S^{(0)} \downarrow S^{(1)} \uparrow S^{(2+)}$, or to the ferromagnetic avalanche $S^{(0)} \downarrow S^{(1)} \downarrow S^{(2-)}$ (Fig. 6b). Therefore, for each transition $S^{(0)} \rightarrow S^{(1)}$, there is one possible ferromagnetic and one antiferromagnetic avalanche when $S^{(1)}$ is not a saturated state. In the case where $S^{(1)}$ is saturated, only the antiferromagnetic option exists. We must check for each of these transitions that $S^{(2)} \neq S^{(0)}$, because an avalanche cannot contain any cycles. We thus obtain a list of avalanches for each base graph, from which we construct a list of candidate graphs by evaluating each combination of avalanches.

An example for one of the $N = 2$ base graphs is shown in Fig. 7. For the example shown, the base graph has four possible $l = 2$ avalanches all associated with a different transition ($00 \uparrow 01 \uparrow 11$, $01 \uparrow 11 \downarrow 10$, $11 \downarrow 10 \downarrow 00$, and $10 \downarrow 00 \uparrow 01$), and therefore $2^4 = 16$ candidate graphs.

Thus far, we have only considered $L = 2$ graphs. We now describe how we construct graphs that have avalanches of length $L = 3$ based on the $L = 2$ graphs.

Suppose that we have a $L = 2$ graph. Similarly to how we can find $l = 2$ avalanches for a base graphs by evaluat-

ing combinations of incoming and outgoing transitions, we can construct possible $l = 3$ avalanches by finding all combinations of $l = 2$ avalanches and outgoing $l = 1$ transitions for the given graph. From the possible $l = 3$ avalanches, we can again construct a list of candidate graphs.

We now ask which $L = 2$ graphs should be used as our new 'base' **AVOID CONFUSION W/ BASE GRAPHS** for the $L = 3$ graphs. It is insufficient to use only the $L = 2$ graphs that are themselves realizable, as it is not guaranteed that a derived $L = 3$ graph will also be impossible. We observe that, when extending a $l = 2$ avalanche to a $l = 3$ avalanche, most of the conditions required for the $l = 2$ avalanche are still included in the conditions for the $l = 3$ avalanche (see SI). The only exception is the final set of relaxation inequalities that describes the stability of the state $S^{(2)}$, which is stable for the $l = 2$ transition but unstable for the $l = 3$ transition. As a result, we can predict that for a $L = 2$ graph of which the base inequalities and the first set of relaxation inequalities are incompatible, any derived graphs with longer avalanches must also be impossible. If, however, it is only the last set of relaxation inequalities that causes the graph to be impossible, then a derived graph of $L = 3$ or higher may still be realizable.

We can thus distinguish three groups of $L = 2$ graphs, as indicated in Fig. 7. The first group of graphs is realizable, and can therefore also be used as a base to construct $L = 3$ graphs. For the second group, the graphs themselves are not realizable, but derived $L = 3$ graphs may be realizable. For the third group, neither the $L = 2$ graphs themselves nor any derived graphs containing longer avalanches are realizable. To construct all possible $L = 3$ graphs, we only need to use graphs from the first two groups.

We now generalize our method. To find all $L = l$ graphs, we construct the graphs for $L = l - 1$, and divide the graphs into the three groups as described. We then construct all $L = l$ candidate graphs for each of the $l = l - 1$ graphs in the first two groups. We thus have an algorithm that constructs all realizable graphs by iterating over the maximum avalanche length L . Our algorithm terminates when none of the graphs belonging to the first and second groups have any possible avalanches of length l .

IV. APPLICATION TO SPECIFIC MODELS

We now illustrate how the base graph method is applied to the specific model where the coupling between hysterons is linear. Linear coupling entails that the switching fields $U_i^{+,-}(S)$ depend on the state S as:

$$U_i^{+,-}(S) = u_i^{+,-} - \sum_{i \neq j} c_{ij} s_j \quad (13)$$

Depending on the system, the allowed coupling coefficients c_{ij} may be restricted. We consider two instances of the linear model for $N = 3$. First, we look at the case

where only one of the coupling coefficients c_{ij} is unequal to zero. Second, we look at a model for hysterons in series, where the coupling is $c_{ij} = c_j$ with all $c_j < 0$.

A. Linear coupling

B. Hysterons in series

V. DISCUSSION

VI. OUTLOOK

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