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Monte Carlo investigation of how interfacial magnetic couplings affect blocking temperature distributions in exchange bias bilayers

G. Lhoutellier,¹ D. Ledue,^{1,a)} R. Patte,¹ and V. Baltz^{2,3,4,b)}

¹Normandie Univ., INSA Rouen, UNIROUEN, CNRS, GPM, 76000 Rouen, France

²Univ. Grenoble Alpes, SPINTEC, F-38000 Grenoble, France

³CNRS, SPINTEC, F-38000 Grenoble, France

⁴CEA, INAC-SPINTEC, F-38000 Grenoble, France

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Exchange bias in ferromagnetic (F)/antiferromagnetic (AF) bilayers is a function of both the bulk properties of the AF layer and the interfacial properties determining the effective interfacial couplings between the F and AF layers. The distinction between bulk and interface can be clearly revealed in blocking temperature distributions, where AF grain volume distribution results in a high-temperature peak while disordered interfacial magnetic phases produce a low-temperature contribution. However, the coupling conditions producing such bimodal blocking temperature distributions remain to be specified. In this article, we use a granular model which accounts for the disordered interfacial phases by considering small magnetic grains (SGs) with weaker anisotropy and coupling with the F grains at the F/AF interface. The SG are included in the AF material. The coupling conditions producing bimodal blocking temperature distributions were determined. Then, using Monte Carlo simulations, these conditions were validated and the effect of interfacial F-SG coupling on distributions was investigated. We next determined how the ratio between F-SG and F-AF couplings could be used to estimate the surface coverage of the disordered interfacial phases from experimental data. *Published by AIP Publishing.* [<http://dx.doi.org/10.1063/1.4967829>]

I. INTRODUCTION

Exchange bias refers to the magnetic exchange interactions between ferromagnetic (F) and antiferromagnetic (AF) materials in a F/AF bilayer. F/AF exchange biased bilayers are used in spintronic devices to set the reference direction determining the spin of conduction electrons.^{1,2} Exchange bias is determined by numerous parameters such as the bulk anisotropies, the bulk and interfacial exchange couplings, and the grain volumes in polycrystalline films.^{1,2} In the AF layer, two categories of magnetic phases are usually considered: bulk phases^{3–6} and interfacial phases with a certain level of surface coverage.^{7–13} These phases have low freezing temperatures due to magnetic frustration produced by defects at the F/AF interface, and are also known as disordered interfacial spins or low freezing temperature spins. These interfacial phases can significantly affect the exchange bias properties. At working temperatures for devices, they no longer contribute to the exchange field.¹⁴

Experimentally, the magnetic phases in the AF layer are accessible through several techniques, for example, they can be deduced from thermoremanent studies down to low temperatures.¹⁰ Although this method is appropriate for isolated AF materials, it does not work for F/AF bilayers because the signal from the F layer screens that of the disordered phases. Another technique, x-ray magnetic dichroism, can measure uncompensated AF spins in magnetic phases.⁹ However, it requires large-scale facilities, and x-ray absorption by the

top layers in a stack may create difficulties. An alternative is to measure blocking temperature distributions to identify magnetic phases in the AF layer.^{5,8} This technique presents no particular technical difficulties, and can be performed, for example, using the stepwise procedure described by Soeya *et al.*⁵ In this procedure, the sample is first field-cooled from above the maximum blocking temperature down to 4 K under a positive field. This cooling causes the magnetic phases of the AF layer in contact with the F layer to align in the positive direction. Then, the sample's temperature is raised to an intermediate annealing temperature T_a , and the sample is subsequently field-cooled under a negative field down to 4 K. This cooling step results in a negative reorientation of the magnetic phases of the AF layer in contact with the F layer, when the blocking temperature is lower than T_a . Finally, a hysteresis loop is measured at 4 K. These steps are repeated for a range of T_a and the variation in the exchange field, H_E , with T_a and its derivative dH_E/dT_a are recorded. Typical experimental results for Co(3 nm)/IrMn(7 nm) bilayers¹⁵ have been reproduced in Fig. 1. Since the increase in the exchange field with T_a is related to the negative repolarization of the magnetic phases of the AF layer in contact with the F layer, the derivative dH_E/dT_a , is linked to the blocking temperature distributions of these magnetic phases.^{5,8} The volume distribution of grains in the AF bulk phases in contact with the F layer gives rise to a high-temperature peak; a low-temperature contribution is also observed, which is attributed to the interfacial phases. The relative contribution from the interfacial phases can be extracted from the relative weights of the high- and low-temperature contributions.¹⁴ However, the impact of the

^{a)}denis.ledue@univ-rouen.fr

^{b)}vincent.baltz@cea.fr

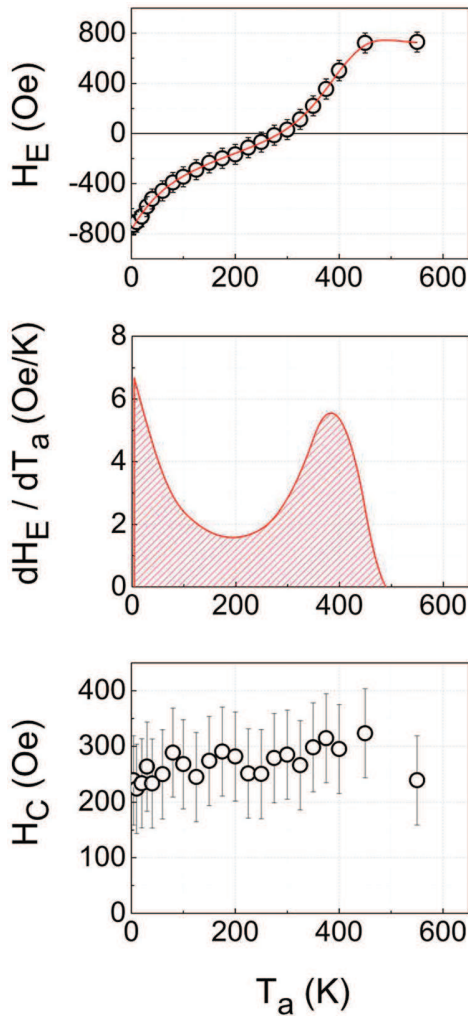


FIG. 1. Experimental dependences of H_E (top, the full lines result from interpolation of the data), the derivative dH_E/dT_a (middle, from the interpolation) and H_C (bottom) on T_a for Co(3 nm)/IrMn(7 nm) bilayers (from Ref. 15).

interfacial couplings on each contribution remains to be determined.

From a theoretical point of view, it is not easy to account for such double-peaked dH_E/dT_a . For example, attempts were made to use a single effective F/AF coupling which slowly decreases as the temperature increases.³ Unfortunately, this model produced a single peak. Recently, a Monte Carlo investigation based on a granular model¹⁶ validated the hypothesis that less stable magnetic regions in the AF layer, located at the F/AF interface, could be responsible for the low-temperature peak of dH_E/dT_a and thus, the bimodal distribution of blocking temperature, experimentally observed.

In the present study, using a granular model, we, first, analytically determined the interfacial coupling conditions required to obtain a double-peaked derivative dH_E/dT_a . These predictions were subsequently validated using Monte Carlo simulations which are very suitable for simulating the procedure described above.⁵ In addition, we discuss the relationship between the derivative dH_E/dT_a and the blocking temperature distribution. We also indicate how our results

may be useful in refining how the relative amount of interfacial magnetic phases can be determined from experimental data.

II. MODEL AND SIMULATION

Our model bilayer (Fig. 2) consisted of one F layer and one AF layer with the same microstructure since we assumed columnar growth. For each layer, the granular microstructure was generated by Voronoi tessellation¹⁷ in two-dimensions and the resulting volume distribution was log-normal. The average volume is $(56.25 \times t) \text{ nm}^3$ where t is the layer thickness. The standard deviation σ is about 0.43 and the median grain size is $(51.28 \times t) \text{ nm}^3$. Smaller grains (SGs) with reduced anisotropy compared to the bulk AF grains were randomly spread over the F/AF interface within the AF (Fig. 2). The surface coverage of these SG was expressed as a percentage of the total surface, x_{SG} . The SG account for the magnetic phases with low freezing temperatures due to magnetic frustration produced by defects at the F/AF interface (e.g., interlayer diffusion and stacking faults). Due to that, the SG are not compositionally similar to the AF but show deteriorated magnetic properties. Since the dimensions of the grains are in the nanometer scale, every grain was assumed to support a single magnetic domain that reverses by uniform rotation. Thus, a unit vector, σ_i , represents the magnetization orientation of each F grain, each SG and the interfacial uncompensated magnetization orientation of each bulk AF grain. Across the interface, each F grain was coupled with one AF grain (J_{F-AF}), or with one SG (J_{F-SG}). In addition, each SG was coupled with one AF grain below (J_{AF-SG}). In agreement with the literature, the AF grains were not linked to each other,^{6,18,19} and nor were the SG. In contrast, the F grains were coupled to each other (J_F) (Fig. 2). Uniaxial anisotropy along a common easy axis (y-axis) in the plane of the layer was considered for all grains, and a linear thermal dependence of the anisotropy constants per unit volume K_i , was implemented.²⁰ Thus, the total energy of the system can be expressed as

$$E = - \sum_{(i,j)} J_{ij} \sigma_i \cdot \sigma_j - \sum_i K_i V_i (\sigma_i \cdot \mathbf{e}_y)^2 - \mu_0 \mathbf{H} \cdot \sum_{i \in F, SG} \mathbf{m}_i,$$

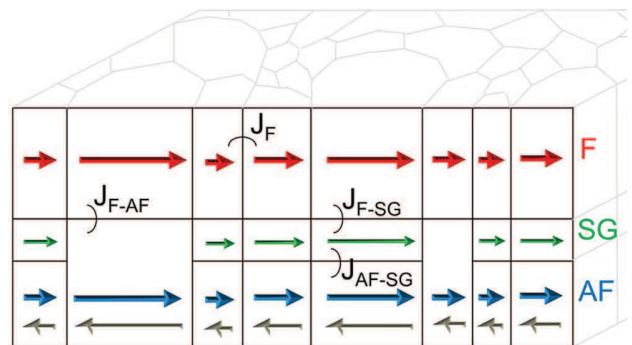


FIG. 2. Schematic representation of the F/AF bilayer. The model considers F and AF grains as well as smaller grains (SG) present at the F/AF interface.

where V_i is the grain volume, \mathbf{m}_i is the magnetic moment of the F grain or SG, and \mathbf{H} is the field applied along the y -axis.

Next, the protocol described by Soeya *et al.*⁵ was simulated using the Monte Carlo method.^{21,22} At each temperature (during the field-cooling) or at each value of the applied field (during the hysteresis loop), n Monte Carlo steps are performed, that is, each grain is visited n times (typically, n is several thousand). At each visit, a magnetization reversal trial with an acceptance rate equal to $\exp(-\Delta E/(k_B T))$, where ΔE is the energy barrier to overcome during reversal, is made.^{16,23} Readers should remember that these energy barriers mainly depend on the anisotropy and also on the local field acting on the grain considered. In the Monte Carlo method, the time unit is one Monte Carlo step. Unfortunately, the physical time δt corresponding to one Monte Carlo step is, in general, unknown. However, in our particular case, it has been found that $\delta t \approx 2\tau_0$ ²⁴ where τ_0 is the prefactor in the expression of the relaxation time $\tau = \tau_0 e^{K_V/k_B T}$. Consequently, the physical time corresponding to n Monte Carlo steps is $n \delta t \approx 2 n \tau_0$, which corresponds to the measuring time t_m .

To compare our results with available experimental results,^{8,14,15} we used 5-nm-thick F and AF layers. Assuming that the interfacial magnetic phases extended over 3–4 atomic planes, we set the thickness of the SG, t_{SG} , to 1 nm,^{25,26} thus, reducing the thickness of the AF grains located below them to 4 nm (see Fig. 2). Since the study was based on Co/IrMn bilayers, the transition temperatures used in the expression of the anisotropy constants²⁰ were $T_C = 1360$ K for the F grains (Co) and $T_N = 690$ K for the AF grains (IrMn).¹ The 0 K anisotropy constant for the AF grains was estimated to be $K_{AF}^0 = 9.8 \times 10^5 \text{ J m}^{-3}$.²⁷ We then took $K_F^0 = 7.2 \times 10^4 \text{ J m}^{-3}$ so that the intrinsic blocking temperature would be 140 K for a F grain of volume equal to the average volume, $\bar{V}_F = (7.5 \times 7.5 \times 5) \text{ nm}^3$. The effective anisotropy of the SG was assumed to be weaker than that of the AF grains; we, therefore, arbitrarily set K_{SG}^0 to $0.3 K_{AF}^0$. As a basis for further discussion, we calculated the intrinsic blocking temperature distribution for the AF grains and SG, that is, the blocking temperatures in the absence of coupling and applied field. The intrinsic blocking temperature $T_{B,int}$ of a grain is defined by $\tau(T_{B,int}) = t_m$, where $\tau = \tau_0 e^{K_V/k_B T}$. The intrinsic blocking temperatures of the AF grains and of the SG can thus be written as

$$T_{B,int} = \frac{KV}{k_B \ln(t_m/\tau_0)},$$

where $K = K_{AF}$ or K_{SG} . Since, in our simulations, $t_m \approx 2 n \tau_0$, we could calculate the intrinsic blocking temperature distribution of the grains from their volume distributions. These distributions for $n = 5.5 \times 10^3$ are plotted in Fig. 3. The intrinsic blocking temperatures of an AF grain of volume equal to the average volume $\bar{V}_{AF} = (7.5 \times 7.5 \times 5) \text{ nm}^3$, and a SG grain of volume equal to its average volume $\bar{V}_{SG} = (7.5 \times 7.5 \times 1) \text{ nm}^3$, will be $T_{B,int}^{AF} = 521.6$ K and $T_{B,int}^{SG} = 108.4$ K, respectively. Below, we will show that even though the intrinsic blocking temperature distribution is bimodal, it is not guaranteed that dH_E/dT_a , obtained from

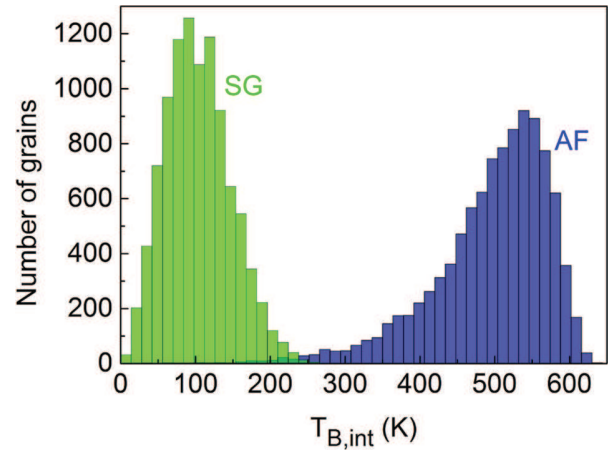


FIG. 3. Intrinsic blocking temperature distributions for the SG and AF grains calculated from the volume distribution simulated for $N = 10^4$ grains.

Ref. 5 procedure, will have a double-peak, since this characteristic depends on the degree of coupling between grains.

The effective interfacial coupling per unit area j_{F-AF} , is simply related to the exchange field in the case of an ideal F/AF interface, that is, without SG. Here, j_{F-AF} was chosen so that the value of H_E at 0 K, when all AF grains are polarized in the same direction, would be equal to 700 Oe, in agreement with previous experimental studies¹⁵ (Fig. 1). We, therefore, took $j_{F-AF} = 5 \times 10^{-22} \text{ J nm}^{-2}$. The actual couplings per unit area, j_{F-SG} and j_{AF-SG} , are unknown. However, j_{F-SG} is assumed to be smaller than j_{F-AF} . In addition, j_{F-SG} and j_{AF-SG} must be chosen such that: (i) the SG should be partially repolarized toward the negative direction when field-cooling is applied (the fraction of SG that will be repolarized depends on T_a and on the field applied \mathbf{H}_{FC}) and (ii) the coercive field should not vary with T_a , that is, with the fraction of repolarized entities in contact with the F layer. This second condition is required for the model to be consistent with the experimental data^{5,8,15} (Fig. 1). Condition (i) is fulfilled if and only if the local field (including the applied field) acting on a SG is negative and smaller in absolute value than the anisotropy field $H_a^{SG} = \frac{2K_{SG}}{\mu_0 M_{SG}}$ (M_{SG} is the magnetization of a SG, $M_{SG} = 164 \text{ kA m}^{-1} = M_{AF}/8$ to take into account magnetic disorder such as interlayer diffusion and stacking faults). Since the local field acting on a SG is the sum of the field due to the F grain in contact $\mathbf{H}_{F-SG} = \frac{J_{F-SG}}{\mu_0 M_{SG} V_i} \boldsymbol{\sigma}_F$ (with $J_{F-SG} > 0$ and $\boldsymbol{\sigma}_F = -\mathbf{e}_y$), the field due to the AF grain in contact $\mathbf{H}_{AF-SG} = \frac{J_{AF-SG}}{\mu_0 M_{SG} V_i} \boldsymbol{\sigma}_{AF}$ (with $J_{AF-SG} > 0$ and $\boldsymbol{\sigma}_{AF} = \mathbf{e}_y$), and \mathbf{H}_{FC} (Fig. 4(a)), this condition can be written as

$$-H_a^{SG} < -\frac{J_{F-SG}}{\mu_0 M_{SG} V_i} + \frac{J_{AF-SG}}{\mu_0 M_{SG} V_i} - |H_{FC}| < 0,$$

with $V_i = S_i t_{SG}$ (S_i is the surface of the SG). The condition can be rewritten as

$$0 < j_{F-SG} - j_{AF-SG} + \mu_0 M_{SG} t_{SG} |H_{FC}| < 2K_{SG}(T) t_{SG}. \quad (1)$$

Condition (ii) requires that the negatively repolarized SG should not switch magnetization after the F grains have

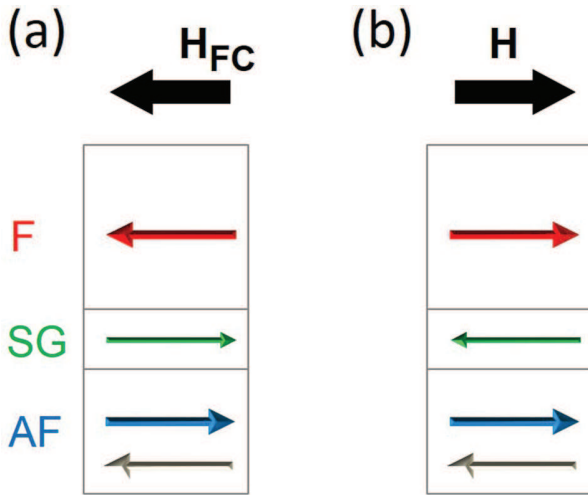


FIG. 4. Schematic representation of the F/AF bilayer during (a) field cooling before negative reorientation of the SG and (b) the hysteresis loop at 4 K after reversal of the magnetization from negative to positive direction in the F layer.

reversed their magnetization from negative to positive orientation during the hysteresis loops (right hand side of the hysteresis loop). Thus, the local field (including the applied field) acting on a SG after the F grains have reversed (Fig. 4(b)) should be smaller than the anisotropy field H_a^{SG} . This condition can be written as

$$0 < \frac{J_{F-SG}}{\mu_0 M_{SG} V_i} + \frac{J_{AF-SG}}{\mu_0 M_{SG} V_i} + H < H_a^{SG},$$

where $H > 0$ is the field applied during the hysteresis loop, and thus

$$0 < j_{F-SG} + j_{AF-SG} + \mu_0 M_{SG} t_{SG} H < 2 K_{SG}(T) t_{SG}. \quad (2)$$

It should be noted that this inequality is a necessary condition, but the coercive field can vary if the F layer fails to reverse by uniform rotation. Inequalities (1) and (2) are summarized in Fig. 5. According to this diagram, only the conditions represented in the green area (lower left corner) will produce double-peaked dH_E/dT_a with a constant coercive field as T_a varies (assuming uniform rotation of the F layer). We remark that increasing the anisotropy constant of the SG is equivalent to decreasing j_{F-SG} and j_{AF-SG} and to shifting upward the blue long-dashed line and to shifting downward the black full and the red short-dashed lines in Fig. 5.

III. RESULTS AND DISCUSSION

To check whether in the conditions indicated, double-peaked or single-peaked dH_E/dT_a will be produced, we considered three pairs of (reduced) couplings ($j_{F-SG}/(K_{SG}^0 t_{SG})$, $j_{AF-SG}/(K_{SG}^0 t_{SG})$): (1.8, 0.1) (blue area in Fig. 5); (0.2, 1.2) (orange area in Fig. 5); and (0.2, 0.1) (green area in Fig. 5).

The simulated variation of H_E vs. T_a was obtained from the simulated hysteresis loops at 4 K (not shown here). Then, we numerically calculated the derivative of its interpolation. The variation in H_E and its derivative were plotted against T_a for $j_{F-SG}/(K_{SG}^0 t_{SG}) = 1.8$ and $j_{AF-SG}/(K_{SG}^0 t_{SG}) = 0.1$, and $j_{F-SG}/$

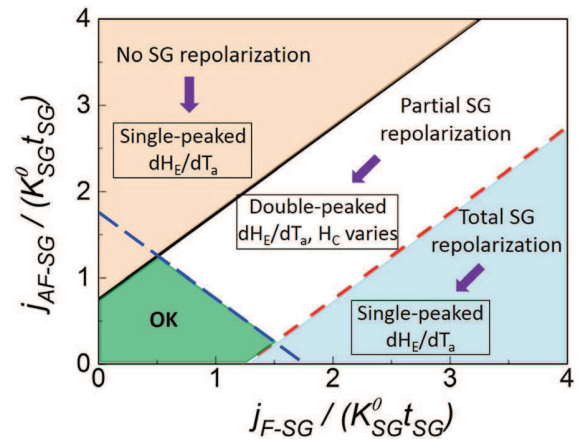


FIG. 5. dH_E/dT_a as a function of the reduced variables $j_{F-SG}/(K_{SG}^0 t_{SG})$ and $j_{AF-SG}/(K_{SG}^0 t_{SG})$ (the black full line corresponds to a local field acting on a SG equal to 0, the increasing red short-dashed line to a negative local field acting on a SG whose absolute value is equal to the anisotropy field H_a^{SG} , after negative repolarization of the F layer during field cooling, and the decreasing blue long-dashed line to a local field acting on a SG equal to the anisotropy field H_a^{SG} , after reversal of the magnetization in the F layer during the hysteresis loop). The green zone (lower left corner) represents the conditions in which dH_E/dT_a will display a double peak.

$(K_{SG}^0 t_{SG}) = 0.2$ and $j_{AF-SG}/(K_{SG}^0 t_{SG}) = 1.2$ (Fig. 6, left and middle columns, respectively). In these figures, as expected, there is no peak at low-temperature, even though SG are present in large amounts ($x_{SG} = 75\%$) at the F/AF interface. However, the reasons for the absence of a low-temperature peak are not the same in the two different cases. Thus, in Fig. 6 (left column), there is no peak at low-temperature because, due the strong j_{F-SG} coupling, all SG are repolarized by the F grains during field-cooling even for $T_a = 4$ K. Therefore, even if there is no visible peak at low-temperature, the exchange field is affected by the SG at room temperature. The absence of a low-temperature peak in Fig. 6 (middle column) is because the SG cannot be repolarized while the AF grains are not repolarized due to the strong j_{AF-SG} coupling in this case. In these conditions, the SG start to be repolarized only from $T_a = 270$ K. The contribution of the SG to the derivative dH_E/dT_a is masked by the contribution of the AF grains, and the exchange field at room temperature is not affected by the SG.

Thus, a double-peaked dH_E/dT_a is expected only for the set $j_{F-SG}/(K_{SG}^0 t_{SG}) = 0.2$ and $j_{AF-SG}/(K_{SG}^0 t_{SG}) = 0.1$. The variation of H_E and its derivative dH_E/dT_a in these conditions is shown in Fig. 6 (right column). In agreement with the predictions from our analytical calculations, two contributions are observed. In addition, the locations of the two peaks are in good agreement with the intrinsic blocking temperature distributions shown in Fig. 3, thus, confirming the validity of the procedure. However, a slight shift of the low-temperature peak toward lower than expected temperatures is observed due to the J_{F-SG} coupling which lowers the energy barrier for negative repolarization of the SG. We will discuss the relative heights of the two peaks below. In addition, the coercive field is independent of T_a as in previous experimental work^{5,8,15} (Fig. 1).

The fraction of SG (x_{SG}) at the interface needs to be extracted from the area of the low-temperature peak of dH_E/dT_a which is the quantity measured experimentally. To see

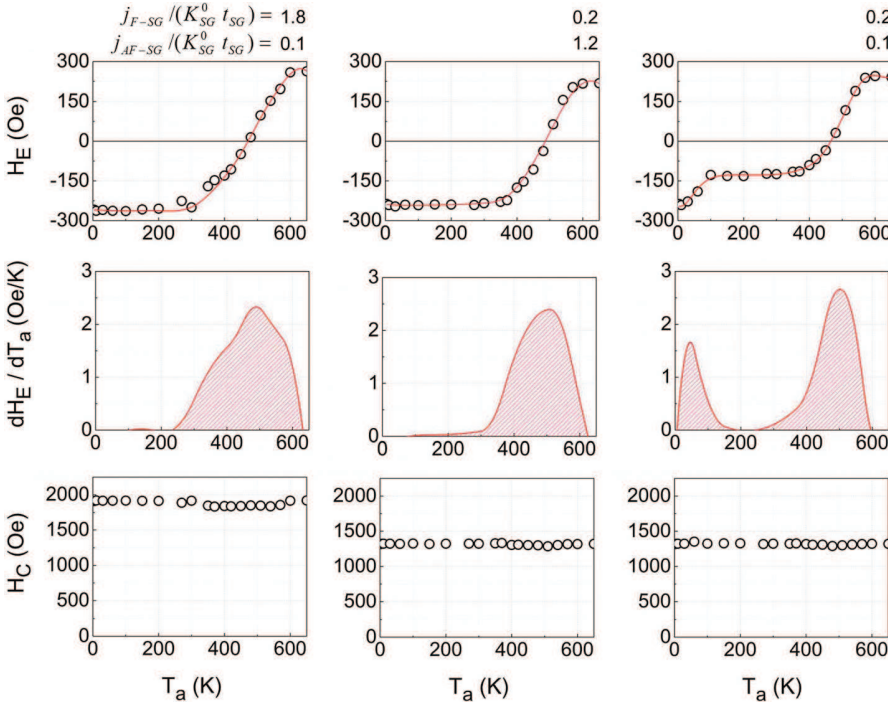


FIG. 6. Variation of H_E vs. T_a , the full line is an interpolation of the data (top row), derivative dH_E/dT_a vs. T_a (middle row, from the data interpolation) and variation of H_C vs. T_a (bottom row) for three pairs of (reduced) couplings ($j_{F-SG}/(K_{SG}^0 t_{SG})$, $j_{AF-SG}/(K_{SG}^0 t_{SG})$) and $x_{SG} = 75\%$.

how this can be done, we investigated the effect of the j_{F-SG} coupling ($j_{F-SG}/(K_{SG}^0 t_{SG}) = 0.1, 0.2,$ and 0.4 , that is, $j_{F-SG}/j_{F-AF} = 0.05, 0.1,$ and 0.2) on the derivative dH_E/dT_a for $x_{SG} = 75\%$ and $j_{AF-SG}/(K_{SG}^0 t_{SG}) = 0.1$ (Fig. 7). From this representation, it is clear that as j_{F-SG} increases, the low-temperature peak of the derivative related to the SG grows, whereas the high-temperature peak (of the derivative) due to the AF grains remains constant. If we define the quantity $\Delta^* = S_L/(S_H + S_L)$, where S_L and S_H are the surfaces of the low- and high-temperature peaks, respectively (Fig. 7), we obtain $\Delta^* = 13\%, 22\%,$ and 35% for $j_{F-SG}/j_{F-AF} = 0.05, 0.1,$ and 0.2 , respectively. It should be noted that, as j_{F-SG} increases, for the blocking temperature distribution, only a small shift toward low temperatures and no growth of the low-temperature peak is expected; the high-temperature peak should be unaltered because the blocking temperatures

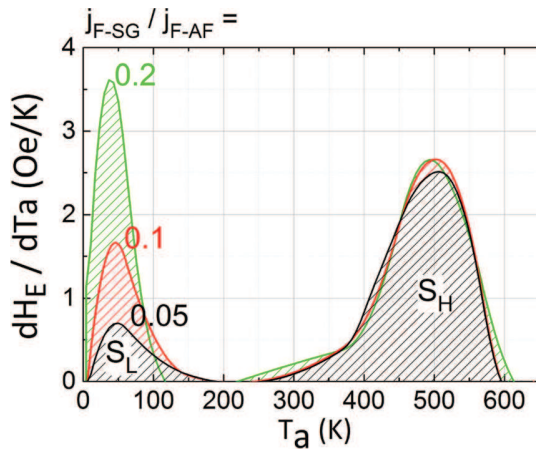


FIG. 7. Derivative dH_E/dT_a calculated from H_E vs. T_a for various intensities of F-SG coupling, $j_{F-SG}/j_{F-AF} = 0.05, 0.1,$ and 0.2 , with $x_{SG} = 75\%$ (S_L and S_H correspond to the surfaces of the low- and high-temperature peaks).

of the AF grains are not affected by j_{F-SG} . The results presented above clearly indicate that the derivative dH_E/dT_a is not proportional to the blocking temperature distribution.

To define how x_{SG} , the quantity of interest, is related to the experimentally measured Δ^* , we calculated the expression of dH_E/dT_a assuming a uniform rotation of the F layer (see Appendix A), and found

$$\frac{dH_E(T_a)}{dT_a} = \frac{2}{\mu_0 M_F t_F} [x_{SG} j_{F-SG} P_{SG}(T_a) + (1 - x_{SG}) j_{F-AF} P_{AF}(T_a)],$$

where M_F is the magnetization of each F grain, t_F is the thickness of the F layer, and P_{SG} and P_{AF} are the normalized blocking temperature distributions of the SG and AF grains, respectively ($\int_0^{+\infty} P_{SG}(T) dT = \int_0^{+\infty} P_{AF}(T) dT = 1$). The surface fraction of the low-temperature contribution can readily be obtained from the above expression of dH_E/dT_a (see Appendix B), thus

$$\Delta^* = [x_{SG} j_{F-SG}/j_{F-AF}] / [1 - x_{SG} + x_{SG} j_{F-SG}/j_{F-AF}].$$

The variation in Δ^* and its relationship with x_{SG} for various values of j_{F-SG}/j_{F-AF} is illustrated in Fig. 8. From this representation, it is clear that Δ^* is lower than x_{SG} assuming $j_{F-SG} < j_{F-AF}$, and this difference increases as j_{F-SG}/j_{F-AF} decreases in agreement with the simulated Δ^* values for $j_{F-SG}/j_{F-AF} = 0.05, 0.1,$ and 0.2 ($x_{SG} = 75\%$) (Fig. 8). Note that $\Delta^* = x_{SG}$ only if $j_{F-SG} = j_{F-AF}$. This was to be expected because in that case dH_E/dT_a is proportional to the blocking temperature distribution.

IV. CONCLUSION

In the framework of a granular model, and using Monte Carlo simulations, we specified some conditions on the

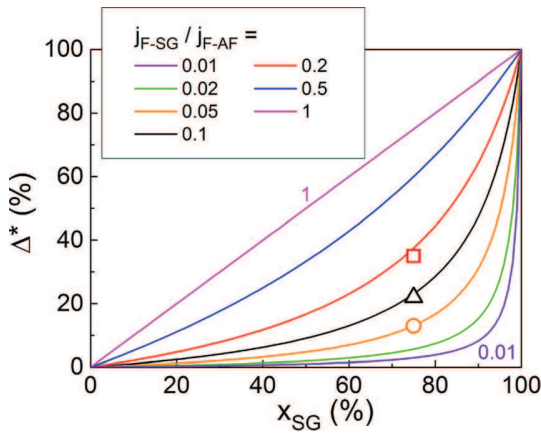


FIG. 8. Calculated variation in the surface fraction Δ^* of the low-temperature peak of dH_E/dT_a with proportion of SG at F/AF interface, x_{SG} , for various j_{F-SG}/j_{F-AF} ratios. The open symbols represent values obtained from Monte Carlo simulations.

coupling intensity between small grains (SG) taking into account less stable phases at a ferromagnetic (F)/antiferromagnetic (AF) interface and the F or AF grains, which explain experimental results indicating a bimodal blocking temperature distribution. In addition, in these conditions, the coercive field is not dependent on the annealing temperature as for experimental data. Importantly, at 0 K these two couplings must be approximately smaller or equal to the anisotropy energy of the SG ($K_{SG}^0 V_{SG}$). The surface coverage of disordered magnetic phases at the interface can, on principle, be extracted from experimental measurements using the protocol presented by Soeya *et al.*,⁵ but our model demonstrated that this calculation requires the ratio between F-SG and F-AF couplings to be known. Alternatively, this ratio can be extracted from experimental data, on the condition that it is possible to vary only the surface coverage of the SG without affecting other parameters.

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APPENDIX A: RELATIONSHIP BETWEEN THE DERIVATIVE dH_E/dT_a AND THE BLOCKING TEMPERATURE DISTRIBUTION

Let us consider the F/AF bilayer in Fig. 2 with a fraction x_{SG} of SG at the interface. Assuming that a fraction r of AF grains in contact with the F layer and a fraction r' of SG are negatively repolarized, the average interfacial coupling per unit surface is

$$\begin{aligned} j_{\text{int}} &= [-r(1-x_{SG}) + (1-r)(1-x_{SG})]j_{F-AF} \\ &\quad + [-r'x_{SG} + (1-r')x_{SG}]j_{F-SG} \\ &= (1-2r)(1-x_{SG})j_{F-AF} + (1-2r')x_{SG}j_{F-SG}. \end{aligned}$$

If the F layer reverses by uniform rotation, the exchange field H_E at 0 K will be given by

$$\mu_0 H_E M_F t_F = (2r-1)(1-x_{SG})j_{F-AF} + (2r'-1)x_{SG}j_{F-SG}, \quad (\text{A1})$$

where M_F is the magnetization of each F grain and t_F is the thickness of the F layer.

The derivative of the exchange field therefore becomes

$$\mu_0 M_F t_F \frac{dH_E}{dT_a} = 2 \left[(1-x_{SG})j_{F-AF} \frac{dr}{dT_a} + x_{SG}j_{F-SG} \frac{dr'}{dT_a} \right]. \quad (\text{A2})$$

The fractions r and r' are related to the normalized blocking temperature distributions for SG and AF grains in contact with the F layer

$$r = \int_0^{T_a} P_{AF}(T_B) dT_B, \quad r' = \int_0^{T_a} P_{SG}(T_B) dT_B,$$

with $\int_0^{+\infty} P_{SG}(T) dT = \int_0^{+\infty} P_{AF}(T) dT = 1$. Consequently, $\frac{dr(T_a)}{dT_a} = P_{AF}(T_a)$ and $\frac{dr'(T_a)}{dT_a} = P_{SG}(T_a)$.

Equation (A2) can therefore be rewritten as

$$\begin{aligned} \frac{dH_E(T_a)}{dT_a} &= \frac{2}{\mu_0 M_F t_F} \left[(1-x_{SG})j_{F-AF} P_{AF}(T_a) \right. \\ &\quad \left. + x_{SG}j_{F-SG} P_{SG}(T_a) \right], \end{aligned}$$

and the derivative is proportional to the blocking temperature distribution $= (1-x_{SG})P_{AF}(T_B) + x_{SG}P_{SG}(T_B)$ only if $j_{F-AF} = j_{F-SG}$.

APPENDIX B: EXPRESSION TO DETERMINE THE SURFACE FRACTION OF THE LOW-TEMPERATURE PEAK

The surface fraction, Δ^* , of the low-temperature peak of dH_E/dT_a is equal to $S_L/(S_H + S_L)$ where $S_L = H_E(T_a = 200 \text{ K}) - H_E(T_a = 4 \text{ K})$ and $S_H + S_L = H_E(T_a = 650 \text{ K}) - H_E(T_a = 4 \text{ K})$. Assuming uniform rotation of the F layer, $H_E(T_a = 200 \text{ K})$ can be obtained from Equation (A1) in Appendix A, with $r=0$ and $r'=1$ (all SG are negatively repolarized, but the AF grains in contact with the F layer are not), $H_E(T_a = 4 \text{ K})$ can be obtained with $r=r'=0$ since all the entities in contact with the F layer will be positively oriented, while $H_E(T_a = 650 \text{ K})$ can be obtained with $r=r'=1$ since all the entities in contact with the F layer will be negatively repolarized (this gives $H_E(T_a = 650 \text{ K}) = -H_E(T_a = 4 \text{ K})$). In this case, Δ^* is given by

$$\Delta^* = \frac{x_{SG}j_{F-SG}/j_{F-AF}}{1-x_{SG} + x_{SG}j_{F-SG}/j_{F-AF}}.$$

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