

# Simulating magnetic hysteresis in the Stoner-Wohlfarth Model using Fortran 90

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The Stoner-Wohlfarth model is a simple model for a magnetic hysteresis, involving a single-domain particle with magnetocrystalline anisotropy in an external magnetic field.

## Goal:

- Correctly set-up the energy functional for the Stoner-Wohlfarth particle
- Implement a reliable energy minimizer
- Calculate the hysteresis and other relevant quantities, such as the energy profiles and the Stoner-Wohlfarth Astroid

The inspiration for this topic was taken from a python simulation of the SW model I did in the course "VU Micromagnetics and Spintronics: Models and Simulation".

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- Ferromagnetic materials form so-called domains, clusters in which many magnetic moments align parallel. Under the influence of an external magnetic field, these domains quickly all orientate themselves in the direction of the field.

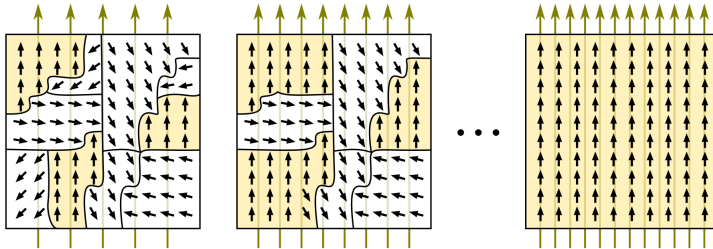


Figure: Magnetic domains in a ferromagnet and their response to an applied field.

Source: Wikimedia commons, 'Growing magnetic domains', 2020. Author: MikeRun,  
<https://commons.wikimedia.org/wiki/File:Growing-magnetic-domains.svg>

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- The particle possesses magnetocrystalline anisotropy along its easy axis, i.e. the magnetization 'likes' to orient itself along this direction, as this minimizes its total energy.
- Despite not being accurate enough to model hysteretic behavior on a large scale, it is a very good model for small magnetic particles used in magnetic storage, magnetic sensors, nanotechnology and biomagnetism. Furthermore, it is an extremely useful model for understanding magnetic anisotropy.

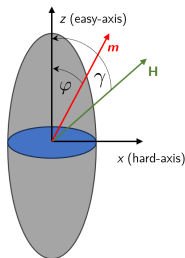


Figure: A sketch of a stoner-Wohlfarth particle.

- The energy per volume unit  $e = E/V$  of the SW particle can be derived as

$$e = e_{\text{aniso}} + e_{\text{ext}},$$

and these energetic contributions (derivation skipped) lead to an energy functional dependent only on the angle  $\varphi$ ,

$$e(\varphi) = k \cdot \sin^2 \varphi - \mu_0 M_s (H_{\parallel} \cos \varphi + H_{\perp} \sin \varphi). \quad (1)$$

- In equation 1,  $H_{\parallel} = H \cos \gamma$  and  $H_{\perp} = H \sin \gamma$  are the components of the external field  $\mathbf{H}$  parallel and perpendicular to the easy axis, respectively.
- Often, one is interested in the critical 'switching field', i.e. the strength of the field  $\mathbf{H}$  required to reorient the magnetization in its direction. If we consider the energy, this switching field is given by the saddle points of the functional, i.e one must find  $\frac{\partial e}{\partial \varphi} = 0$  and  $\frac{\partial^2 e}{\partial \varphi^2} = 0$ .

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- The hysteresis itself can be simulated by minimizing the energy as a function of  $\varphi$  for a fixed angle  $\gamma$  and a range of field strengths  $H \rightarrow$  the obtained  $\varphi$  can be used to determine the magnetization orientation at the energetic minimum (magnetization in field direction  $\mathbf{m}_H = \cos \varphi$ ).

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**What is a good choice for a minimizer?** Secant method or derivative-based method?

We go for an iterative derivative-based minimizer, precisely, the conjugate gradient method.



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- The exact algorithm used is the Fletcher-Reeves algorithm, which calculates mutually conjugate search directions  $\delta_j$  in each iteration,

$$\delta_{j+1} = \begin{cases} -\nabla f(x_j), & \text{if } j = 0 \\ -\nabla f(x_j) + x_j \delta_j, & \text{if } j = 1, 2, \dots, n-1. \end{cases} \quad (2)$$

In the above equation,

$$x_j = x_{j-1} + \alpha_j \delta_j, \quad (3)$$

where  $\alpha_j$  is the  $j$ th step length.  $x_j$  in Eq. 2 is given by,

$$\alpha_j = \frac{\|\nabla f(x_j)\|^2}{\|\nabla f(x_{j-1})\|^2}. \quad (4)$$