

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".

- Magnetic materials consist of a large number of so-called 'elementary magnets', i.e. magnetic dipole moments of each atom which interact with each other and external magnetic fields.
- In electrodynamics of continua, we define the magnetization field \mathbf{M} as

$$\mathbf{M} = \frac{d\mathbf{m}}{dV},$$

where \mathbf{m} is a magnetic moment and dV is the infinitesimal volume element.

- A magnetic hysteresis curve describes how the magnetic moments of a material behave in response to an external magnetic field, in general, this depends on the nature of a magnetic material (dia-, para-, and ferromagnetic) and other properties.
- 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.

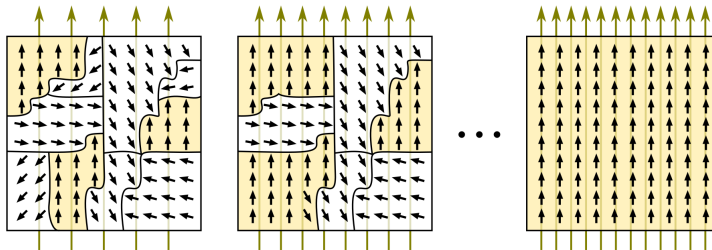


Figure: Magnetic domains in a ferromagnet and their response to an applied field.

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

- Describing the hysteresis of a ferromagnetic material is theoretically very complex and involves various energetic contributions, and a simulation requires complex numerical treatment to accurately resolve all effects
→ *micromagnetic* theory and simulations
- We can turn to a simpler model, which is the Stoner-Wohlfarth model, describing a single-domain ellipsoid magnetic particle.
- 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.

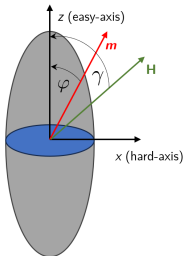


Figure: A sketch of a stoner-Wohlfarth particle.

- 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 φ ,

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

- 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 $\frac{\partial^2 e}{\partial \varphi^2} = 0$.
- The hysteresis itself can be simulated by minimizing the energy as a function of α for a fixed angle γ and a range of field strengths $H \rightarrow$ the obtained α can be used to determine the magnetization orientation at the energetic minimum (magnetization in field direction $\mathbf{m}_H = \cos \alpha$).

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.

- Maybe a bit 'overkill'/unnecessary in a one-dimensional problem, but provides very stable results, which is not always the case with simple gradient descent. Also requires less input arguments than secant methods, which makes code more simple
- The exact algorithm used is the Fletcher-Reeves algorithm, which calculates mutually conjugate search directions δ_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 α_j is the j th step length. x_j in Eq. 2 is given by,

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