Debugging time evolution

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1 Checking periodically perturbed free energy

To check that the configuration force matches with the energy, we check the energy for a uniform configuration first, and then for a small perturbation. The latter check gives an estimate for the configuration force by the functional Taylor series expansion:

$$F(Q + \delta Q) = F(Q) + \int_{\Omega} \left(\frac{\delta F}{\delta Q} \Big|_{Q} : \delta Q \right) dV + \mathcal{O}\left(|\delta Q|^{2} \right)$$
 (1)

with

$$\frac{\delta F}{\delta Q} = \frac{\partial f}{\partial Q} - \nabla \cdot \frac{\partial f}{\partial (\nabla Q)} \tag{2}$$

Since this holds for an arbitrary functional F, and also for an arbitrary (small) perturbation δQ , it's probably true that:

$$f(Q + \delta Q) = f(Q) + \left(\frac{\partial f}{\partial Q} - \nabla \cdot \frac{\partial f}{\partial (\nabla Q)}\right) : \delta Q + \mathcal{O}\left(|\delta Q|^2\right)$$
(3)

for each component of the free energy f. For the example we test, we take Q to be:

$$Q = S_0 \begin{bmatrix} \frac{2}{3} & 0 & 0\\ 0 & -\frac{1}{3} & 0\\ 0 & 0 & -\frac{1}{3} \end{bmatrix}$$
 (4)

with

$$\delta Q = S_0 \begin{bmatrix} 0 & \epsilon \sin kx & 0\\ \epsilon \sin kx & 0 & 0\\ 0 & 0 & 0 \end{bmatrix}$$
 (5)

This corresponds (to first order) to a perturbation of the director so that the total configuration $(Q + \delta Q)$ is a uniform-S configuration with director $\mathbf{n} = (1, \epsilon \sin kx, 0)$.

Here, the nondimensional free energy is given by:

$$F = \int_{\Omega} \left(-\frac{1}{2}\alpha Q : Q + \left(\log 4\pi - \log Z + \Lambda : \left(Q + \frac{1}{3}I \right) \right) + \frac{1}{2} \left| \nabla Q \right|^2 \right) dV \tag{6}$$

Then we can calculate F(Q) partially analytically and partially numerically. Of course the elastic part is zero, and we calculate:

$$-\frac{1}{2}\alpha Q: Q = -S_0^2 \frac{\alpha}{3} = -1.2153600266666669$$
 (7)

with $S_0 = 0.6751$ and $\alpha = 8.0$. Further, we can calculate numerically (see calc_lambda program with generate_periodic_Q_tensors.py) Λ and Z for $S_0 = 0.6751$:

Z = 3.87017170996747

$$\Lambda = \begin{bmatrix}
3.6005952163635766 & 0 & 0 \\
0 & -1.8002976081817883 & 0 \\
0 & 0 & -1.8002976081817883
\end{bmatrix}$$
(8)

Then we may calculate F by multiplying this energy density by the size of the domain. The uniform configuration check is then just to compare that number with the energy output from the simulation for a uniform configuration with the same S_0 and α values and domain size. Calculating the energy density from the mean-field interaction gives:

$$f_{\text{entropy}} = 3.60848720197831 \tag{9}$$

Running the simulation and calculating the free energy over the domain gives:

$$F_{\text{mean field}} = -11.99497833$$
 (10)

and calculating out the energy density from the entropy term gives:

$$F_{\text{entropy}} = 35.61419603$$
 (11)

Comparing these values to the energy density multiplied by $(3.1415926 \times 3.1415926)$ (the domain size) we get similar results.

We may also consider a configuration with a different S_0 -value, just to be sure. Running a similar calculation with $S_0 = 0.5$, we get the following values for Λ and Z:

Z = 1.8088523960817302

$$\Lambda = \begin{bmatrix}
2.323990879382531 & 0 & 0 \\
0 & -1.1619954396912653 & 0 \\
0 & 0 & 0
\end{bmatrix}$$
(12)

Then the entropy and mean field part of the free energy density is given by:

$$f_{\text{entropy}} = 3.100327077782174$$

 $f_{\text{mean field}} = -\frac{2}{3}$ (13)

Both of these correspond with the energy densities written to the vtu file.

The configuration force, which is exactly the functional derivative of the free energy, is given by:

$$\frac{\delta F}{\delta Q} = -\alpha Q + \Lambda - \nabla^2 Q \tag{14}$$

We need to evaluate this with the unperturbed Q-tensor – since this is just a uniform configuration, the elastic term will be zero. However, if we explicitly calculate it we find that:

$$\left. \frac{\delta F}{\delta Q} \right|_{Q} : \delta Q \tag{15}$$

Hence, we actually can't get any information out of this particular perturbation.