Ulmer_Project

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Ulmer_Project

Analyzing the influence of non-neutral contact angle on the Cahn-Hilliard model

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Contents

for example, a square and a circle.

condition

Abstract

This project examines the effect of ~~a simple~~ Neumann boundary ~~approach~~ on a finite difference solver for the Cahn-Hilliard equation. It presents the results of the proposed method on different domains (~~a square and a circular example~~). There it is apparent, that the approach is able to ~~effect~~ the contact angle of the interface on the boundary. Additionally, this project gives a rudimentary technique to calculate this angle.

affect

Single coupled

1 Introduction

(CH)

This project thesis builds upon the work in our bachelor thesis, by introducing a simple boundary condition approach to a variation of the solver used therein. In Chapter 2 we introduce the Cahn-Hilliard equation in the formulation that we use for this project. This project used a ~~two dimensional~~ second order version of this CH equation rather than the usual ~~1D~~ 4th order one, to simplify the numerical implementation. The solver itself builds upon a finite difference discretization of this equation. In Chapter 4, together with the discrete domains, on which we run our numerical solver, we introduce a ~~Jacoby~~ iteration to solve the linear system derived from the aforementioned discretization. The numerical solver in this thesis is GPU accelerated, and the discretization we chose to base our solver on is capable to calculate on all domains as long as a characteristic function is given. Therefore, we introduce two domains, on which we present our findings. The primary goal of this work is then the boundary condition approach in Chapter 3. Conceptionally the boundary condition we introduce simply consists of a constant value added in the linear system to all equations corresponding to grid-cells on the boundary. The actual implementation is capable of doing this for arbitrary domains. We present the results of this method on two Domains, a square one in 5 and a circular domain in 6 where we show the phase field for different boundary conditions which manifest in a variable contact angle of the interface on the boundary mimicking the behavior of hydrophobic/hydrophilic material. While we are unable to provide explicit formulae in relation to the ~~constant~~, in Chapter 7 we provide numerical insight in this relationship, and a table with precomputed values.

Jacobi

Chapte 5

Chapter 6

contact angle

2 Fundamentals

This work concerns itself with boundary conditions on the Cahn-Hilliard equation. The Cahn-Hilliard (CH) equation is a fourth order partial differential equation (PDE) used to describe phase seperation in binary mixtures. It models how a mixture of two components (e.g., two liquids or alloys) evolves over time to separate into distinct regions with different concentrations of each component. To achive this it provides a phase-field ϕ which is used for an implicit representation of the interface between both phases. The Cahn-Hilliard equation, in the formulation we use here, is derived from the **Ginzburg-Landau** energy (2.1), an example on how this is done is given by [3].

$$E^{\text{bulk}}[\phi] = \int_\Omega \frac{\varepsilon^2}{2}|\nabla\phi|^2 + W(\phi)\,dx, \tag{2.1}$$

There they introduce a chemical potential μ derived as derivative of the **Ginzburg-Landau** energy.

$$\mu = \frac{\delta E_{bulk}(\phi)}{\delta\phi} = -\varepsilon^2\Delta\phi + W'(\phi), \tag{2.2}$$

Where $W(\phi)$ in the energy, is a double well potential. In our case we orient us at the work of[1], where they use

$$W(\phi) = \frac{(1-\phi^2)^2}{4}. \tag{2.3}$$

the Cahn-Hilliard equation in this thesis is then given as

$$\begin{aligned} \partial_t\phi(x,t) &= \Delta\mu \\ \mu &= -\varepsilon^2\Delta\phi + W'(\phi). \end{aligned} \tag{2.4}$$

One thing to note is, that this way of writing the CH equation presents a second order, ~~two-dimensional~~ system, rather than the one ~~dimensional~~ fourth order system (2.5) often given.

coupled

$$\partial_t \phi(\vec{x}, t) = \Delta(-\varepsilon^2 \Delta\phi + W'(\phi)) \tag{2.5}$$

This choice is deliberate, and aligns with the numerical implementation.

2.1 Notation

This project solves the CH equation on a regular rectangular grid with grid-size h. The computational domain, is therefore discretized as

$$\vec{x}_{ij} := \frac{i}{h} * e_1 + \frac{j}{h} e_2 \qquad \frac{j}{h} * e_2 \tag{2.6}$$

where $i, j \in [0, \dots N]$ and N is chosen arbitrarily, such that the resulting rectangle $[0, Nh] \times [0, Nh]$ acts as bounding box of the domain Ω. For our implementation we use $N = 256$ as it gives a good compromise between resolution and compute time. We denote a discrete version of the domain Ω_d where

$$\Omega_d := \{x_{ij} | x_{ij} \in \Omega\} \tag{2.7}$$

On this discrete domain our solver calculates solutions for discrete fields

$$\phi_{ij}^n : \Omega_d \times \{0, \dots\} \to \mathbb{R}, \tag{2.8}$$
$$\phi_{ij} := \phi(\vec{x}_{ij}) \qquad \vec{x}_{ij} \in \Omega_d \tag{2.9}$$
$$\mu_{ij}^n : \Omega_d \times \{0, \dots\} \to \mathbb{R}, \tag{2.10}$$
$$\mu_{ij} := \mu(\vec{x}_{ij}) \tag{2.11}$$

We use the following differential quotients for field f_{ij}:

$$D_x f_{i+\frac{1}{2}j} = \frac{f_{i+1j} - f_{ij}}{h} \qquad D_y f_{ij+\frac{1}{2}} = \frac{f_{ij+1} - f_{ij}}{h} \tag{2.12}$$

And define a discrete gradient as.

$$\nabla_d f_{ij} = (D_x f_{i+1j},\ D_y f_{ij+1}) \tag{2.13}$$

And

$$\Delta_d f_{ij} = \nabla_d \cdot \nabla_d f_{ij} \tag{2.14}$$

See [2] Our solver implements the ansatz proposed by the authors [1]. write author name or abc et al.

$$
\begin{aligned}
\frac{\phi_{ij}^{n+1} - \phi_{ij}^{n}}{\Delta t} &= \nabla_d \cdot (G_{ij} \nabla_d \mu_{ij}^{n+\frac{1}{2}}) \\
\mu_{ij}^{n+\frac{1}{2}} &= 2\phi_{ij}^{n+1} - \varepsilon^2 \nabla_d \cdot (G_{ij} \nabla_d \phi_{ij}^{n+1}) + W'(\phi_{ij}^{n}) - 2\phi_{ij}^{n}
\end{aligned}
\tag{2.15}
$$

This approach provides a semi implicit time discretization ~~were~~ where linear terms are evaluated implicitly and the nonlinear double well potential is evaluated explicitly.

3 Boundary adaptation

The solver from [1], that we use as reference guaranties no flux boundary conditions at a discrete level by setting $\nabla\phi_{ij} = 0$ for $\phi_{ij} \in \partial\Omega_d$ this is done by multiplying with the characteristic function of Ω_d c

$$G_{ij} = \begin{cases} 1\,, x_{ij} \in \Omega \\ 0\,, x_{ij} \notin \Omega \end{cases} \tag{3.1}$$

To accommodate different boundary conditions, we modify $\nabla_d \cdot (G_{ij}\nabla_d\phi_{ij})$ with a constant term C on grid points next to the boundary. To do this, we introduce a boundary field B_{ij} that we add to μ_{ij}. We determine the value of B_{ij} using a central difference scheme on G

$$B_{ij} = \max\left(|G_{i+\frac{1}{2}j} - G_{i-\frac{1}{2}j}|, |G_{ij+\frac{1}{2}} - G_{ij-\frac{1}{2}}|\right) * C \tag{3.2}$$

For example consider ~~We present an example in case of~~ a 32x32 domain with $C = 1$ of the boundary fields **B** for a square domain 3.1 Figure 3.1 and an example on a circular domain in 3.2. Figure 3.2 In this Project p we use the following adaptation of the discretization from [1].

$$\begin{aligned} \frac{\phi_{ij}^{n+1} - \phi_{ij}n}{\Delta t} &= \nabla_d \cdot (G_{ij}\nabla_d\mu_{ij}^{n+\frac{1}{2}}) \\ \mu_{ij}^{n+\frac{1}{2}} &= 2\phi_{ij}^{n+1} - \varepsilon^2\nabla_d \cdot (G_{ij}\nabla_d\phi_{ij}^{n+1}) + B_{ij} + W'(\phi_{ij}^n) - 2\phi_{ij}^n \end{aligned} \tag{3.3}$$

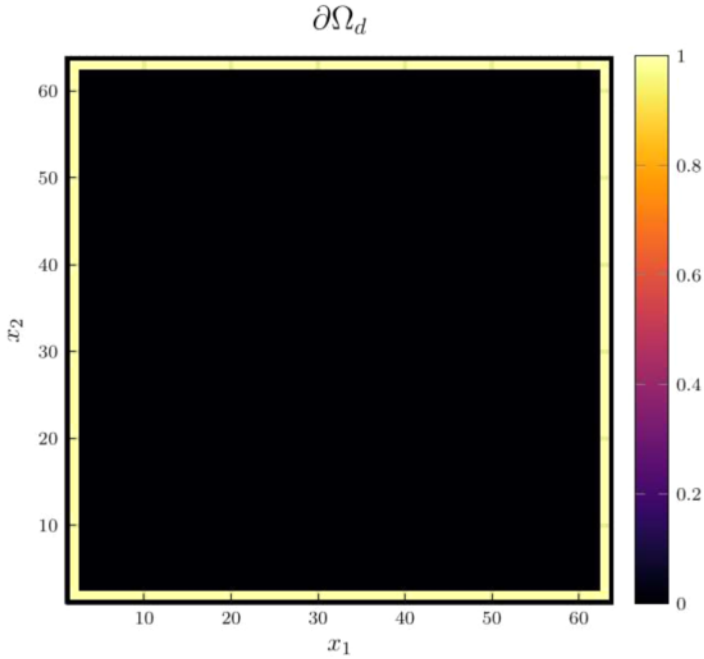


Figure 3.1: Visualization of all grid-cells adjacent to the boundary $\partial\Omega_d$ of a square domain

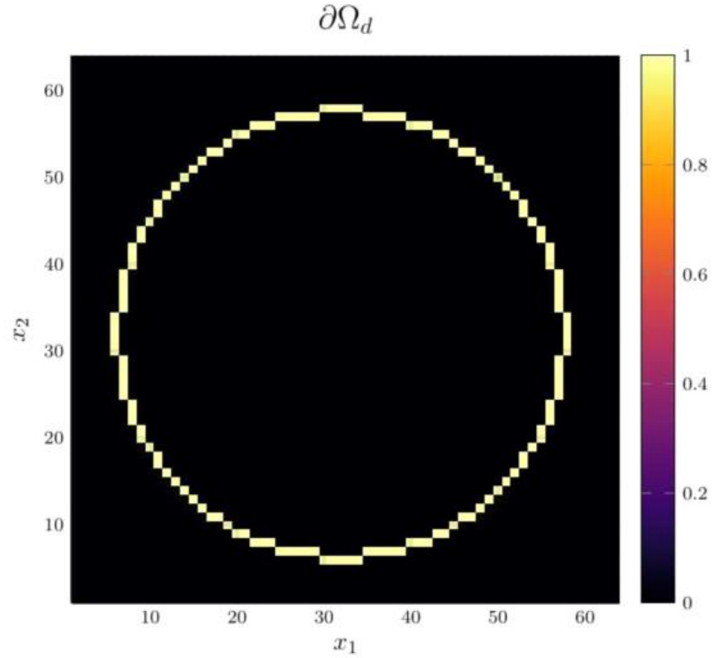


Figure 3.2: Visualization of all grid-cells adjacent to the boundary $\partial\Omega_d$ of a circular domain

4 Numerical solver

Contrary to the solver proposed in [2] we do not use a multi-grid Gauss-Seidel ~~Solver~~ s solver to solve the linear system, instead we use a ~~Jacoby~~ Jacobi solver, as this will eventually assist in parallelizing the computation. Similar to [2] we linearize (3.3) to

$$
\begin{aligned}
\frac{\phi_{ij}^{n+1}}{\Delta t} - \nabla_d \cdot (G_{ij} \nabla_d \mu_{ij}^{n+\frac{1}{2}}) &= \frac{\phi_{ij}^{n}}{\Delta t} \\
\mu_{ij}^{n+\frac{1}{2}} - 2\phi_{ij}^{n+1} + \varepsilon^2 \nabla_d \cdot (G_{ij} \nabla_d \phi_{ij}^{n+1}) + B_{ij} &= 2\phi_{ij}^{n} - W'(\phi_{ij}^{n})
\end{aligned}
\tag{4.1}
$$

Which

One may note, that after rearanging some terms leads to a linear system with a right-hand side ~~wich~~ is exclusively ~~dependant~~ on the previous time step. We use Jacobi's method to solve the resulting linear system given above, and the corresponding element wise representation of the same is given in the following.

dependent

Provided the mth ~~Jacoby~~ Jacobi iteration has been computed, the $m+1th$ iteration is computed by solving

$$
\begin{aligned}
\frac{\phi_{ij}^{n+1,m+1}}{\Delta t} - \nabla_d \cdot (G_{ij} \nabla_d \mu_{ij}^{n+\frac{1}{2},m+\frac{1}{2}}) &= \frac{\phi_{ij}^{n}}{\Delta t} \\
\mu_{ij}^{n+\frac{1}{2},m} - 2\phi_{ij}^{n+1,m} + \varepsilon^2 \nabla_d \cdot (G_{ij} \nabla_d \phi_{ij}^{n+1,m+\frac{1}{2}}) + B_{ij} &= 2\phi_{ij}^{n} - W'(\phi_{ij}^{n})
\end{aligned}
\tag{4.2}
$$

For $\phi_{ij}^{n+1,m+1}, \mu_{ij}^{n+\frac{1}{2},m+1}$, where $\nabla_d \cdot (G_{ij} \nabla_d \mu_{ij}^{n+\frac{1}{2},m+\frac{1}{2}})$ and $\nabla_d \cdot (G_{ij} \nabla_d \phi_{ij}^{n+1,m+\frac{1}{2}})$. can be computed as follows. Use the results from the previous ~~Jacoby~~ Jacobi step for values off the center. e.g.

$$
\begin{aligned}
\nabla_d \cdot (G_{ij} \nabla_d \phi_{ij}^{n+1,m+\frac{1}{2}}) = &\frac{1}{h^2} (G_{i+\frac{1}{2}j} \phi_{i+1j}^{n+1,m} + G_{i-\frac{1}{2}j} \phi_{i-1j}^{n+1,m} \\
&+ \quad G_{ij+\frac{1}{2}} \phi_{ij+1}^{n+1,m} + G_{ij-\frac{1}{2}} \phi_{ij-1}^{n+1,m}) \\
&- \left(G_{i+\frac{1}{2}j} + G_{i-\frac{1}{2}j} + G_{ij+\frac{1}{2}} + G_{ij-\frac{1}{2}}\right) \phi_{ij}^{n+1,m+1}
\end{aligned}
\tag{4.3}
$$

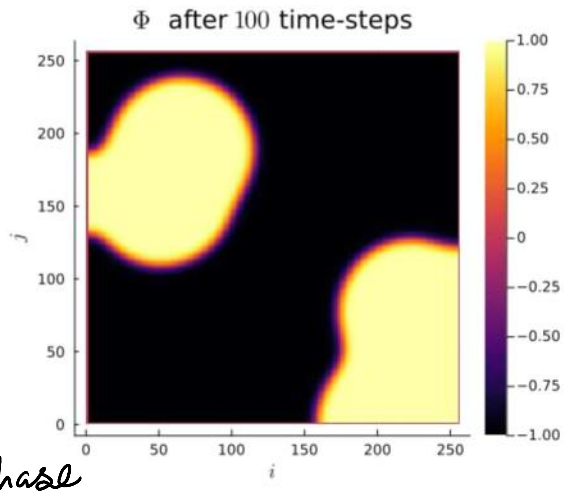
Our implementation is done in julia to transmit the solution for each element in parallel on the GPU. In the following we described the complete implementation of the Jacobi's iteration.

provide some reference for julia

```julia
@kernel function jacoby!(
    Φ,
    M,
    @Const(Ξ),
    @Const(Ψ),
    @Const(h),
    @Const(ε),
    @Const(Δt),
    @Const(iterations)
)
    I   = @index(Global, Cartesian)
    Id  = oneunit(I)
    Ids = CartesianIndices(M)
    Ix = CartesianIndex(1, 0)
    Iy = CartesianIndex(0, 1)
    if I in (Ids[begin]+Id:Ids[end]-Id)
        g = G(2 * I + Ix, Ids) + G(2 * I + Iy, Ids) + G(2 * I - Ix, Ids) + G(2 * I - Iy,
        ↪  Ids)
        a1 = 1/Δt
        a2 = -1* ε^2/h^2 * g  - 2
        b1 = 1/h^2 * g
        b2 = 1
        for _ = 1:iterations

            Σμ = G(2 * I + Ix, Ids) * M[I+Ix] + G(2 * I + Iy, Ids) * M[I+Iy] + G(2 * I -
            ↪  Ix, Ids) * M[I-Ix] + G(2 * I - Iy, Ids) * M[I-Iy]

            Σϕ = G(2 * I + Ix, Ids) * Φ[I+Ix] + G(2 * I + Iy, Ids) * Φ[I+Iy] +G(2 * I -
            ↪  Ix, Ids) * Φ[I-Ix] +G(2 * I - Iy, Ids) * Φ[I-Iy]

            c1 = Ξ[I] + 1/h^2   * Σμ
            c2 = Ψ[I] - ε^2/h^2 * Σϕ

            # stupid matrix solve
            @inline Φ[I] = (c1*b2 - c2*b1) / (a1*b2 - a2*b1)
            @inline M[I] = (a1*c2 - a2*c1) / (a1*b2 - a2*b1)
            #
            @synchronize()
        end

    end
end
```

5 Numerical evaluation

We set constant values for $\mathrm{B_{ij}}$ on the boundary to begin with our evaluations. One may note that $\mathrm{C} = 0$ is equivalent to the no-flux condition of the original solver introduced in the Bachelor thesis. Now, as a preliminary verification step we set $\mathrm{C} = 0$ as our first choice. Consequently, for $C = 0$, the interface lies orthogonal on the boundary (see Fig. 5.1), which we expect for a CH solver with no-flux boundary conditions. For $B_{ij} \in \{-1, 1\}$ we observed behavior connected to hydrophobic / hydrophilic substances on the boundary, where $B_{ij} = 1$ resulted in the one phase pearling off the boundary, while the other seemed attracted. These certainly leads to the apparent contact angles of 180° and 0° respectively. Using $B_{ij} = -1$ results in the opposite behavior.



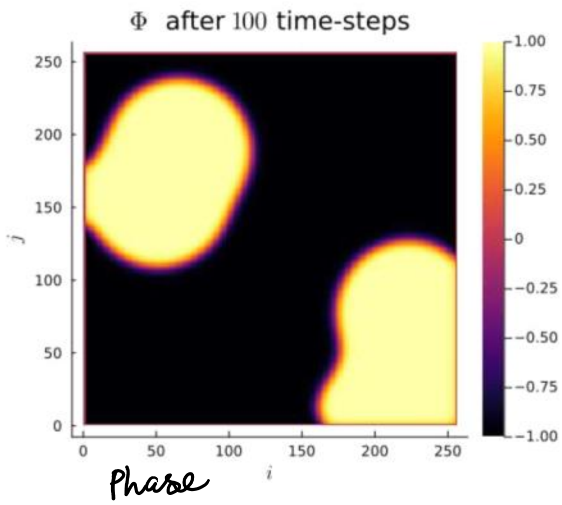
Phase

Figure 5.1: ~~phase~~-field ϕ after 100 time-steps with $C = 0$ emmulating no-flux boundary.

Fig. 5.2

We show, that our solver is stable for values $C \neq 0$. In 5.2 we employ a constant value of $C = 1$ and observe the phase corresponding to $\phi = 1$ puling away from the

boundary. The contact angle between phase 1 and the boundary approaches 180°
i.e. the interface runs parallel to the boundary.



Phase

Figure 5.2: ~~phase~~-field ϕ after 100 time steps with $C = 1$

Fig. 5.3 ,and consequently

the

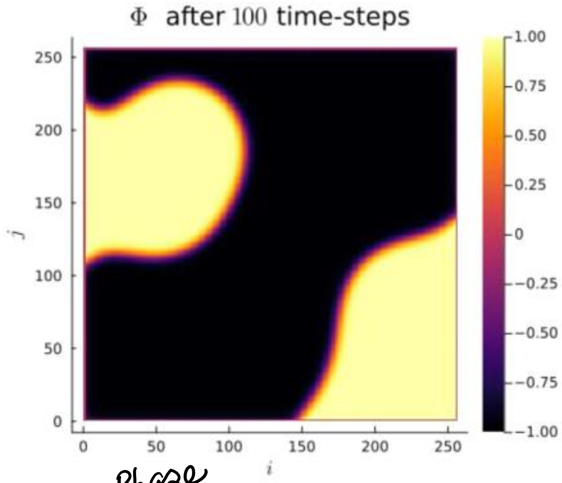
In 5.3 we try the reverse situation. ~~And~~ we observe corresponding behavior. When

the

using a value of $C = -1$ we observe opposite behavior relative to the case in Fig. 5.2. Where the contact angle on the boundary lies at 0°, the interface runs parallel to the boundary again.

The most interesting behavior are noted for values between $(-1, 1)$, where we observe the contact angle of the interface at the boundary changes from parallel 0° to parallel 180°.

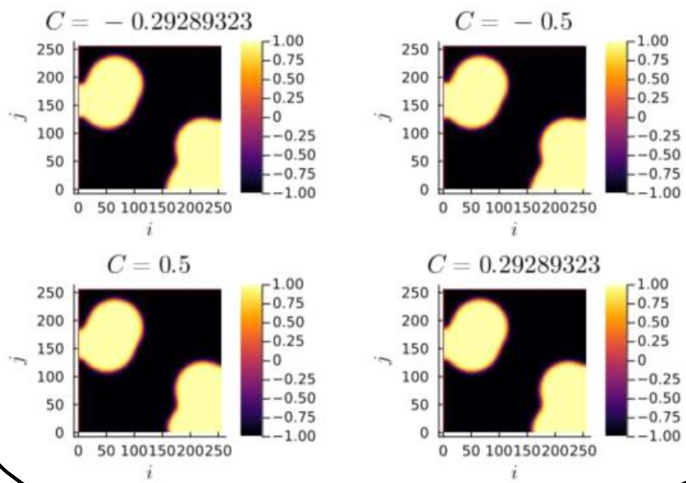
Refer the corresponding figures in the text. Mention that we show the contact angle variation related to different C and h values.



Phase

Figure 5.3: ~~phase~~-field ϕ after 100 time-steps with $C = -1$

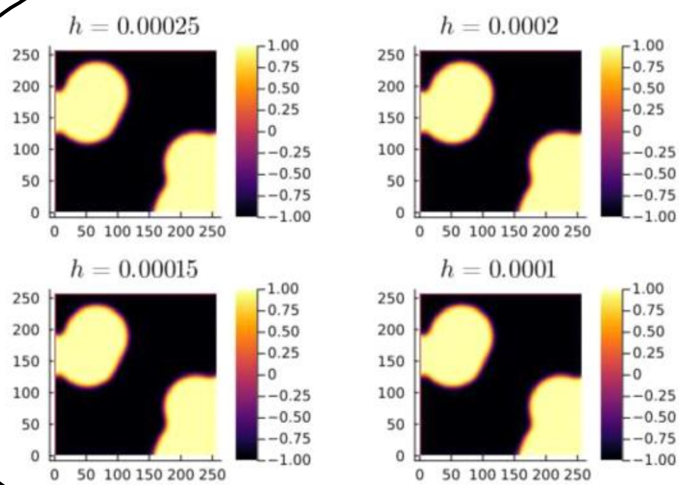
Add figure caption




```
include("src/solvers.jl")
θ = -5f-1
n = 100
arr = _init()
```

```
d = domain(get_backend(arr) , 256 , size(arr))
d(arr)
h = 25e-5
solution = solve(arr , n , θ=θ)
h1 = heatmap(Array(solution) , aspect_ratio=:equal , clims=(-1,1),
↪   lims=(0,size(solution,1)), widen=1.06 , title=L"h=%$h")
h = 20e-5
solution = solve(arr , n , θ=θ)
h2 = heatmap(Array(solution) , aspect_ratio=:equal , clims=(-1,1),
↪   lims=(0,size(solution,1)), widen=1.06 , title=L"h=%$h")
h = 15e-5
solution = solve(arr , n , θ=θ)
h3 = heatmap(Array(solution) , aspect_ratio=:equal , clims=(-1,1),
↪   lims=(0,size(solution,1)), widen=1.06 , title=L"h=%$h")
h = 10e-5
solution = solve(arr , n , θ=θ)
h4 = heatmap(Array(solution) , aspect_ratio=:equal , clims=(-1,1),
↪   lims=(0,size(solution,1)), widen=1.06 , title=L"h=%$h")
plot(h1,h2,h3,h4)
```

Add figure caption



6 Numerical evaluation on a circle

The original solver presented in [1] was able to solve the CH equation on arbitrary domains. Since the addition of our boundary function depends solely on the characteristic function of the discrete domain, we are able to use our approach on different domains ~~Domains~~, by providing a different characteristic function. We present the results Fig. 6.1 of which in this chapter. To show the behavior of the CH solver in 6.1, we first employ no-flux boundary conditions on a circular domain. We observe the interface perpendicular on the boundary, as we expect.

```
W′ (generic function with 1 method)
```

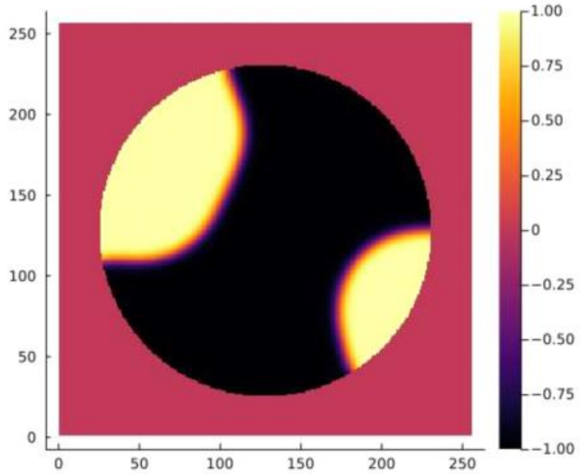
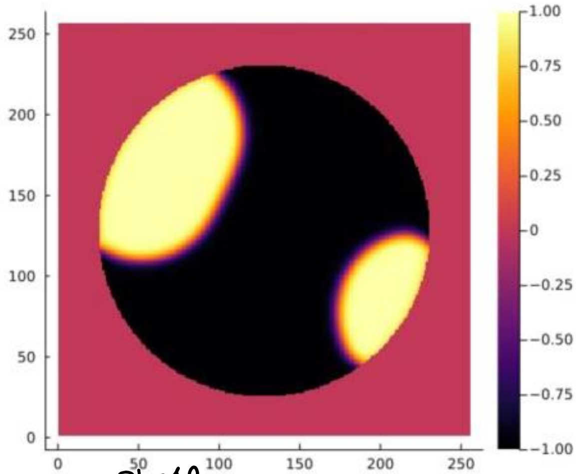


Figure 6.1: ϕ after 100 time steps on a circular domain with no-flux boundary-conditions after 100 time steps on a circular domain with no-flux

Fig 6.2 Fig 5.2

The results we observe in 6.2 are similar to the results on a square domain in 5.2. The contact angle is 180° i.e. the interface does not touch the boundary and runs parallel to it.



Phase

Figure 6.2: ~~phase~~-field ϕ after 100 time-steps with $C = 1$

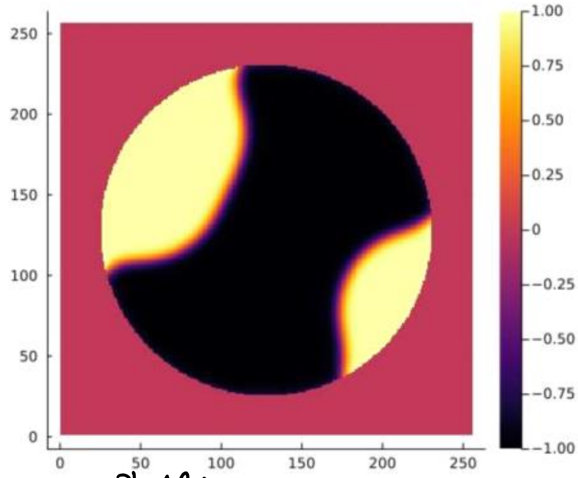
Fig. 5.3 Fig. 6.3

The results for $C = -1$ in 6.3 on the circular domain, are similar to the results in 5.3 on the square domain as well, where the interface touches the boundary and runs parallel with a contact angle of 0°.

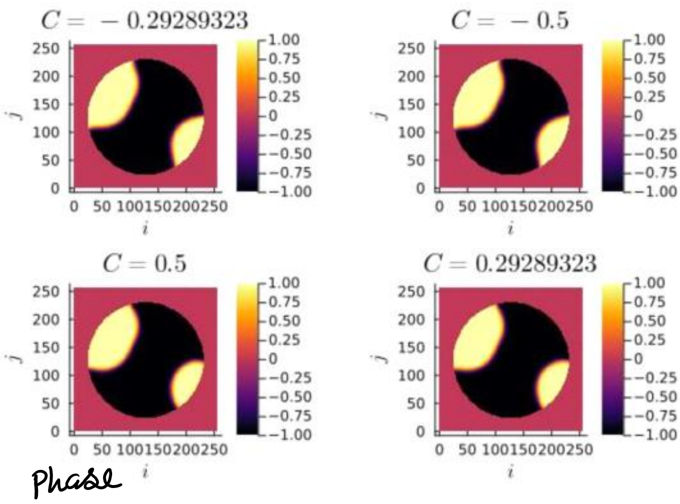
Fig. 6.4

When evaluating intermediate contact angles in 6.4, the results are similar to the square domain again, however, especially for shallow angles, we observe some artifacts of one phase appearing in places where previously was none. We observe similar behavior on square domains only in the corners, i.e. points where the boundary has high curvature.

When using random initial phase-fields, the results look the comparable to the square domain, and exhibit the for the CH equation expected behavior, whereas time goes on, the many small parts coalesce into larger parts.



Phase
Figure 6.3: ~~phase~~-field ϕ after 100 time-steps with $C = -1$



Phase
Figure 6.4: ~~phase~~-field ϕ after 500 time-steps with $C \in \{-1 + \frac{\sqrt{2}}{2}, -0.5, 0.5, 1 - \frac{\sqrt{2}}{2}\}$ on a circular domain.

7 Angle

In previous experiments we noted that the angle of the interface changes with different input parameters. While we do not have a mathematical derivation of this relation, we aim to provide numerical insight in this chapter. We calculate this angle using the gradient of the phase-field $\nabla\phi_{ij}$ and the normal of our domains' boundary.

$$\frac{\nabla_d\phi_{ij}\cdot\mathbf{n}_{ij}}{\|\nabla_d\phi_{ij}\|} = \cos(\theta) \quad \text{where} \quad \phi_{ij}\in\partial\Omega_d \tag{7.1}$$

For a single point $\vec{x}_{ij}$ on the interface and near the boundary. Since we need a finite difference to evaluate 7.1, we do not select a point directly on the boundary and since we need a point on the interface, where $\nabla\phi_{ij}$ is large, we calculate the angle at

Fig. 7.1

$$P_{ij} = arg\max_{\vec{x}_{ij}}\nabla\phi_{ij} \quad \text{where} \quad \phi_{ij}\in\partial\Omega \tag{7.2}$$

```
angle (generic function with 1 method)
```

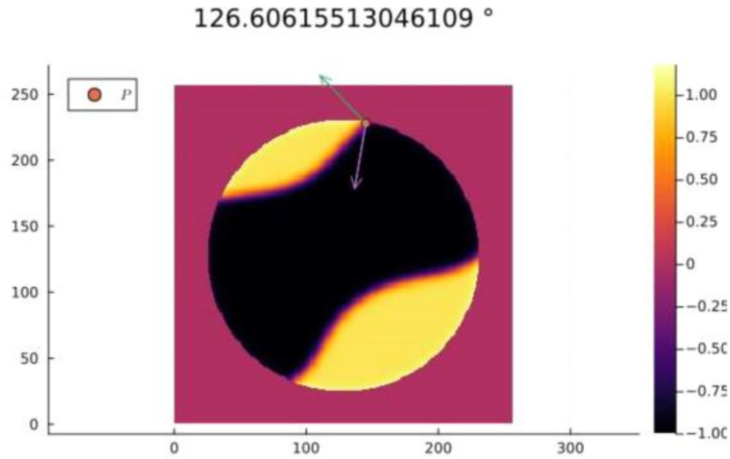
7.1 CIRCLE

The normal of the circular domain in our second example is

$$\mathbf{n}_{ij} := \mathbf{n}(\vec{x}_{ij}) = \frac{\vec{c}-\vec{x}_{ij}}{\|\vec{c}-\vec{x}_{ij}\|} \tag{7.3}$$

Fig. 7.1

Where $\vec{c}$ is the center of the domain. In 7.1 we present the results of a calculated angle, together with the normals and the point it is calculated from.



value

Table 7.1: ~~value~~ for θ and corresponding angle α after 200 time-steps

-0.1	173.49096591056502
-0.095	173.10715739345923
-0.09	172.18364087939332
-0.085	171.54740091859054
-0.08	171.3054040677464
-0.075	171.1455632002332
-0.07	171.02869693204397
-0.065	170.3901810227686
-0.06	170.0449796355949
-0.055	173.27274052589075
-0.05	170.3373892767722
-0.045	168.11953739721892
-0.04	167.41386769034298
-0.035	166.62088559081457
-0.03	164.9014365935728
-0.025	162.8061312020723
-0.02	159.92337650959868
-0.015	155.82320048245077
-0.01	147.4707481361878

-0.005	129.77836444929315
0.0	91.28977210940522
0.005	47.27538237804684
0.01	26.60911004838421
0.015	6.306468865037136
0.02	11.495581754132852
0.025	8.059259459078769
0.03	2.997826637980469
0.035	2.442790881259583
0.04	2.314200756133827
0.045	1.883610279597664
0.05	1.3567468712125557
0.055	0.8024311153759808
0.06	0.5869880299417852
0.065	0.4356076759230446
0.07	0.32719257485287145
0.075	0.03099970458170946
0.08	0.37685133141547533
0.085	0.4151229191583983
0.09	0.7049376111739059
0.095	0.8671639875701463
0.1	1.0282690721714873

7.2 SQUARE

The little

~~the~~ normal vector on a square domain is a ~~litte~~ bit more complicated than the ~~circle~~ normal. For ~~the following~~ we use the normal

for the circle this case

$$\mathbf{n}_{ij} = \mathbf{n}(\vec{x}_{ij}) = \max(\vec{c} - \vec{x}_{ij}) e_{\arg\max_{i,j}(\vec{c}-\vec{x}_{ij})} \tag{7.4}$$

The results are presented in 7.2 Fig. 7.2

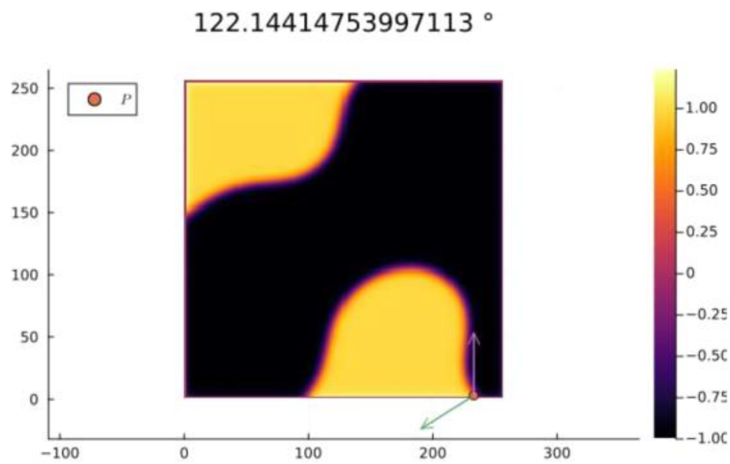


Table 7.2: ~~value~~ Value for θ and corresponding angle α after 200 time-steps

-0.1	177.95311184304808
-0.099	109.40440982717261
-0.098	113.75621160591214
-0.097	105.41178782179448
-0.096	102.88914136385036
-0.095	100.11964229844625
-0.094	103.81071062215845
-0.093	95.69795700045161
-0.092	95.67336816077285
-0.091	99.40375428329858
-0.09	106.41888954576633
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8 Summary and outlook

In this project we examined a numerical model of the CH equation, with simple Neumann boundary conditions. We introduced a simplified version of the solver used in [2] and derived from [1]. Which due to GPU acceleration is significantly faster on our available hardware. We have shown a simple Neumann boundary approach that runs stable on both tested domains. The approach introduced by us is able to freely affect the angle of the phase interface on the boundary, The results of which we have shown on a circular and square domain. We introduced a rudimentary method to calculate the contact angle programmatically however we acknowledge that the results are unreliable. Further research would require a more consistent approach. One such method may be a filter that averages the angle calculation over more than one point, another should be a consistent selection of the point of interest, as the current approach cannot guaranty that the same (or a similar) point is selected if the input parameters change slightly.

Further research may concern itself, with the following topics. First and foremost, we observed inconsistent behavior when changing the hyperparameters ε and grid-size h. However, the methods we used for evaluation were inconsistent at best. Due to the aforementioned unpredictability in the angle calculation the resulting data series was erratic and no trends where apparent. Further research would require investigation of those effects on the boundary. Additionally, in our bachelor thesis, which served as ~~prelimminary~~ work to this project, we investigated an analytical relaxation. The solver used therein for the relaxed problem is compatible with the boundary approach introduced ~~herin~~. Initial tests with the solver for the relaxed system where ~~promissin~~, and further research may investigate those results.

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9 References

Since we have Bibliography
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