

DELAYED HOPF BIFURCATIONS IN REACTION–DIFFUSION SYSTEMS IN TWO SPACE DIMENSIONS

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Abstract

For multi-scale differential equations (or fast–slow equations), one often encounters problems in which a key system parameter slowly passes through a bifurcation. In this article, we show that a pair of prototypical reaction–diffusion equations in two space dimensions can exhibit delayed Hopf bifurcations. Solutions that approach attracting/stable states before the instantaneous Hopf point stay near these states for long, spatially dependent times after these states have become repelling/unstable. We use the complex Ginzburg–Landau equation and the Brusselator models as prototypes. We show that there exist two-dimensional spatio-temporal buffer surfaces and memory surfaces in the three-dimensional space-time. We derive asymptotic formulas for them for the complex Ginzburg–Landau equation and show numerically that they exist also for the Brusselator model. At each point in the domain, these surfaces determine how long the delay in the loss of stability lasts, that is, to leading order when the spatially dependent onset of the post-Hopf oscillations occurs. Also, the onset of the oscillations in these partial differential equations is a hard onset.

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1. Introduction

Many systems of multi-scale differential equations undergo dynamic Hopf bifurcations in which a parameter slowly changes and a stable state becomes unstable due to the

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slow, generic passage of a pair of eigenvalues through the imaginary axis. When the governing equations in the multi-scale (or fast–slow) systems are analytic ordinary differential equations (ODEs), it has been known for over 50 years that there are delays in the onset of the instabilities [4, 20, 35–39, 44–47]. Solutions that have approached the stable (attracting) states before the instantaneous Hopf bifurcations remain near those states long after they have become unstable (repelling) in the dynamic Hopf bifurcations. This phenomenon is known as delayed Hopf bifurcation (DHB).

A primary dynamic manifestation of DHB is that the onset of the post-Hopf oscillations is a hard onset, with the systems transitioning rapidly to large-scale oscillations when they leave neighbourhoods of the unstable states. This contrasts with the gradual square root growth of the oscillation amplitudes in generic classical Hopf bifurcations.

Applications of DHBs in ODE models have been studied in many areas of science and engineering: chemistry [8, 28, 29, 45], electrical circuits [19, 22, 49], electrocardiac models [30], fluid mechanics and geophysics [2, 14, 21, 32], mechanical oscillators [11, 38], neuroscience [4, 5, 9, 10, 24, 42, 43, 46] and economics [18], among others. For each of these models, the long delay between the instantaneous bifurcation and the onset of oscillations, which is $O(1/\varepsilon)$ in the fast time (where ε is the small parameter measuring the time-scale separation), plays an important role in the system dynamics.

Recently, it has been shown that the DHB phenomenon occurs not only in analytic ODEs but also naturally in one-dimensional, multi-scale partial differential equations (PDEs) of reaction–diffusion type [17, 25]. The phenomenon is governed by a super-critical Hopf bifurcation occurring in the family of fast subsystems parametrized by a slowly varying parameter. Examples studied in [17, 25] include the complex Ginzburg–Landau equation, the FitzHugh–Nagumo PDE, and the Hodgkin–Huxley PDE. Additional theory for problems with spectral gaps is given in [3].

In this paper, we show that the phenomenon of DHB also occurs in a pair of multi-scale reaction–diffusion PDEs in two space dimensions. In particular, we study DHBs in the complex Ginzburg–Landau (CGL) PDE in two dimensions and in the Brusselator model in two dimensions. Both of these equations have Hopf bifurcations in which an attracting quasi-steady state (QSS) becomes a repelling QSS, and we consider the problems in which the bifurcation parameters slowly and generically pass through the Hopf points.

First, we use asymptotic analysis on the linearized CGL PDE in two dimensions to show that solutions that approach the attracting QSS before the instantaneous Hopf bifurcation remain close to it even after it has become repelling in the Hopf bifurcation. From the asymptotic analysis, we find that the delays between the time of the instantaneous Hopf bifurcation and the onset of the post-Hopf oscillations are long ($O(1/\varepsilon)$ in the fast time) and generally spatially dependent.

Next, we quantify the duration of these delays. We define spatio-temporal buffer surfaces and spatio-temporal memory surfaces in the three-dimensional space-time of the CGL PDE. These are the surfaces along which the amplitudes of the particular and

homogeneous solutions of the linearized CGL PDE, respectively, cross a threshold, and cause the solution of the full PDE to diverge from the repelling QSS. We derive asymptotic formulas for both surfaces, and we determine how they depend on the initial data and on the properties of the source terms in the CGL PDE. We find that, at each point (x, y) in the domain, numerically computed solutions of the fully nonlinear CGL PDE with initial data given before the Hopf point stay close to the repelling QSS until they reach the spatio-temporal memory surface or the spatio-temporal buffer surface; whichever occurs first at that point. Hence, the linear terms appear to drive the phenomena of DHB in these PDEs, just as they do in the analytic ODEs.

We use the label “buffer” for the spatio-temporal buffer surfaces, because no matter how far in the distant past the data is given (that is, how large negative μ_0 is) the solution must leave the neighbourhood of the repelling QSS at the spatio-temporal buffer surface since the attracting QSS also leaves the neighbourhood then. In addition, we use the label “memory” for the spatio-temporal memory surface, because although the memory of the initial data fades quickly after the initial time due to the rapid, exponential contraction toward the attracting QSS, the memory of the data is not lost. On the contrary, the component of the solution given by the initial data re-emerges in a spatially dependent manner when the slowly varying parameter reaches the memory surface to leading order. This terminology is consistent with that commonly used for DHBs in ODEs.

Finally, we present numerical simulations showing that DHB also occurs in the Brusselator PDE model in two dimensions. Since its creation [41], the Brusselator has served as a prototypical model in pattern formation and chemical oscillation theory (see, for example, [15]). Here, we consider the case when the main bifurcation parameter slowly passes through the Hopf point. As with the CGL PDE, solutions of the Brusselator PDE that approach the attracting QSS before the instantaneous Hopf bifurcation remain near the QSS for long times after it has become repelling. We find that there is a substantial—generally point dependent—delay before the onset of oscillations occurs.

The asymptotic and numerical results that we present here for these prototypical pattern-forming systems in two dimensions extend our earlier work on DHBs in PDEs in one dimension [17, 25]. In particular, in [17] we defined buffer curves and memory curves for the linearized CGL PDE in one dimension and derived asymptotic formulas for their spatio-temporal dependence. We showed that key terms in the particular solution are exponentially small up until the spatio-temporal buffer curve, and that precisely along it these terms become $O(1)$. Similarly, the homogeneous solution is exponentially small up until the memory curve and large after that. Overall, for the CGL PDE in one dimension, we showed that there is a competition between these exponentially small terms: at each point in the one-dimensional domain the term that ceases being small first determines the length of the delay to leading order at that point, and hence also when the hard onset of the (post-Hopf) oscillations occurs. The memory and buffer surfaces defined in this article play a similar role for the CGL PDE in two dimensions.

We recall that classical Hopf bifurcations are ubiquitous in PDEs and spatially extended systems in two dimensions. General results are presented, for example, in [48]. Some specific examples include electrodeposition models [31], bulk-surface reaction–diffusion [40] and nematic liquid crystals [12, 13], in addition to the CGL and Brusselator models studied here.

This paper is organized as follows. In Section 2, we present the main asymptotic analysis to show that DHB occurs in the linearized CGL equation with a slowly varying parameter in two space dimensions. In Section 3, we derive the asymptotic formulas for the spatio-temporal buffer surface and the memory surface for the CGL PDE. In Section 4, we present the results of numerical simulations for the fully nonlinear CGL equation that complement the theoretical predictions (of the previous two sections) for the spatio-temporal dependence of the delayed Hopf bifurcations. Section 5 features the DHB results obtained from numerical simulations of the Brusselator model in two dimensions. Conclusions are presented in Section 6. Some asymptotic calculations are presented in Appendix A.

2. Analysis of DHB in the two-dimensional CGL equation

In two space dimensions, the CGL equation with a source term and a slowly varying parameter is given by

$$\begin{aligned} A_t &= (\mu + i\omega_0)A - (1 + i\alpha)|A|^2A + \sqrt{\varepsilon}I_a(x, y) + \varepsilon d\Delta A, \\ \mu_t &= \varepsilon. \end{aligned} \quad (2.1)$$

Here, $(x, y) \in \mathbb{R}^2$, $\Delta = \partial^2/\partial x^2 + \partial^2/\partial y^2$, $t \geq 0$, $A = A(x, y, t)$ is complex valued, and $0 < \varepsilon \ll 1$ is a small parameter. The linear growth rate $\mu = \mu(t)$ is real for the main phenomena we study. The other system parameters satisfy that $\omega_0 > 0$, α is real and d may be complex valued ($d = d_R + id_I$) with $d_R > 0$, and these parameters are all independent of ε . The source term $I_a(x, y)$ breaks the symmetry $A \rightarrow Ae^{i\theta}$ for any real θ of the CGL equation and is taken to be bounded and positive, with uniformly bounded derivatives. The initial data at $\mu(0) = \mu_0 < 0$ is $A(x, y, 0) = A_0(x, y)$, and taken to be bounded and continuous for all (x, y) . We refer to [1] for the classical CGL PDE.

2.1. The attracting and repelling QSSs The PDE (2.1) has an attracting QSS for all $\mu < -\delta$, where $\delta > 0$ is small and $O(1)$ with respect to ε , and solutions approach it at an exponential rate. Similarly, it has a repelling QSS for all $\mu > \delta$, from which solutions diverge at an exponential rate. These QSS are given by

$$\begin{aligned} A_{\text{QSS}}(x, y) &= -\sqrt{\varepsilon} \frac{I_a(x, y)}{\mu + i\omega_0} \\ &\quad + \varepsilon^{3/2} \left(\frac{I_a(x, y) + d(\mu + i\omega_0)\Delta I_a(x, y)}{(\mu + i\omega_0)^3} - \frac{(1 + i\alpha)I_a^3(x, y)}{(\mu + i\omega_0)^2(\mu^2 + \omega_0^2)} \right) \\ &\quad + O(\varepsilon^{5/2}). \end{aligned} \quad (2.2)$$

Here, the $O(\varepsilon^{5/2})$ terms depend on (x, y) and μ .

2.2. Solution of the linearized equation In this section, we consider $\mu_0 < 0$ and solve the linearized equation,

$$\begin{aligned} A_t &= (\mu + i\omega_0)A + \sqrt{\varepsilon}I_a(x, y) + \varepsilon d\Delta A, \\ \mu_t &= \varepsilon. \end{aligned}$$

This equation may be simplified using an integrating factor. Let $B(x, y, \mu) = A(x, y, \mu)e^{-(\mu+i\omega_0)^2/2\varepsilon}$. The equation for B is an inhomogeneous heat equation,

$$\sqrt{\varepsilon}B_\mu = \sqrt{\varepsilon}d\Delta B + I_a(x, y)e^{-(\mu+i\omega_0)^2/2\varepsilon}. \quad (2.3)$$

The solution of (2.3) with initial data $B(x, y, \mu_0) = A_0(x, y)e^{-(\mu_0+i\omega_0)^2/2\varepsilon}$ is obtained by superimposing the homogeneous solution $B_h(x, y, \mu)$ and the particular solution $B_p(x, y, \mu)$. For $\mu > \mu_0$, we find

$$B_h(x, y, \mu) = \frac{e^{-(\mu_0+i\omega_0)^2/2\varepsilon}}{4\pi d(\mu - \mu_0)} \int_{\mathbf{R}^2} \exp\left[\frac{-(x-x')^2 - (y-y')^2}{4d(\mu - \mu_0)}\right] A_0(x', y') dx' dy'.$$

Here, we have used the fundamental solution of the heat equation in two dimensions,

$$\Phi(x, y, \mu) = \frac{1}{4\pi d\mu} e^{-(x^2+y^2)/4d\mu}.$$

Also, for $\mu > \mu_0$,

$$\begin{aligned} B_p(x, y, \mu) &= \frac{1}{\sqrt{\varepsilon}} \int_{\mu_0}^{\mu} g(x, y, \mu - \tilde{\mu}) e^{-(\tilde{\mu}+i\omega_0)^2/2\varepsilon} d\tilde{\mu}, \\ g(x, y, \mu) &= \frac{1}{4d\pi\mu} \int_{\mathbf{R}^2} \exp\left[\frac{-(x-x')^2 - (y-y')^2}{4d\mu}\right] I_a(x', y') dx' dy'. \end{aligned}$$

Transforming back to the original dependent variable using the integrating factor, we find that, for $\mu > \mu_0$, the homogeneous solution is

$$A_h(x, y, \mu) = \frac{e^{(\mu^2 - \mu_0^2 + 2i\omega_0(\mu - \mu_0))/2\varepsilon}}{4\pi d(\mu - \mu_0)} \int_{\mathbf{R}^2} e^{-(x-x')^2 - (y-y')^2/(4d(\mu - \mu_0))} A_0(x', y') dx' dy'. \quad (2.4)$$

Furthermore, for $\mu > \mu_0$, the particular solution is

$$\begin{aligned} A_p(x, y, \mu) &= \frac{e^{(\mu+i\omega_0)^2/2\varepsilon}}{\sqrt{\varepsilon}} \int_{\mu_0}^{\mu} g(x, y, \mu - \tilde{\mu}) e^{-(\tilde{\mu}+i\omega_0)^2/2\varepsilon} d\tilde{\mu}, \\ g(x, y, \mu) &= \frac{1}{4\pi d\mu} \int_{\mathbf{R}^2} \exp\left[\frac{-(x-x')^2 - (y-y')^2}{4d\mu}\right] I_a(x', y') dx' dy'. \end{aligned}$$

This completes the derivation of the solution of the linearized equation. We calculate A_h for several different types of initial data $A_0(x, y)$ in Section 3.1, and we calculate A_p for several different types of source terms in Section 3.2. Also, we observe here that it will be useful to distinguish between initial data $A_0(x, y)$ given for $\mu_0 \leq -\omega_0$ and for $-\omega_0 < \mu_0 < 0$.

2.3. Solutions stay near the repelling QSS at least until $\mu = \omega_0$ In this section, we consider solutions with $\mu_0 \leq -\omega_0$. We show that not only do the solutions of (2.1) with $\mu_0 \leq -\omega_0$ remain close to the attracting QSS until the time of the instantaneous Hopf bifurcation at $\mu = 0$, but after the parameter crosses the instantaneous Hopf bifurcation they remain close to the repelling QSS as well, at least until the time $\mu = +\omega_0$ at all points (x, y) for the functions $I_a(x, y)$ we consider.

We consider complex values of μ in a horizontal strip with mid-line on the real axis and of sufficient height (at least $3\omega_0$). This enables the use of classical methods of stationary phase and steepest descents (see, for example, [6, 26, 34]) to calculate $A_p(x, y, \mu)$. We find, for $\delta \leq \mu \leq \omega_0$,

$$\begin{aligned} A_p(x, y, \mu) = & -\sqrt{\varepsilon} \frac{I_a(x, y)}{\mu + i\omega_0} + \varepsilon^{3/2} \left(\frac{I_a(x, y) + d(\mu + i\omega_0)\Delta I_a(x, y)}{(\mu + i\omega_0)^3} \right) \\ & + O\left(\frac{\varepsilon^{5/2}}{(\mu + i\omega_0)^5}\right) \\ & + (\sqrt{2\pi}g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}))e^{(\mu + i\omega_0)^2/2\varepsilon}. \end{aligned} \quad (2.5)$$

The calculation is presented in Appendix A.

The first and second terms are precisely the leading-order terms in the expansion of the repelling QSS for the linear CGL (compare with (2.2), where the QSSs are given for the cubic CGL equation). The third (remainder) term contains the higher-order terms in the asymptotic expansion of the repelling QSS, and continued integration by parts will yield them. The fourth term is exponentially small for $\mu \in [\delta, \omega_0 - K\varepsilon^r]$, for some $K > 0$ and any $0 < r < 1$. It is a classic Stokes-type term. This term is not in the expansion of the repelling QSS (on $\mu > \delta$) to all orders. Rather, it is beyond all orders, $O(e^{-\omega_0^2/2\varepsilon})$, arising naturally from tracking solutions on (and near) the attracting QSS along a contour over the saddle point in the complex μ plane and into the regime of $\text{Re}(\mu) > 0$. It is a measure of the exponentially small distance between the attracting and repelling QSS at $\mu = 0$.

Overall, formula (2.5) shows that, at all points (x, y) , the solutions of the linear CGL equation with Gevrey regular data $A_0(x, y)$ given at $\mu_0 \leq -\omega_0$ remain near the repelling QSS at least until $\mu = \omega_0$ to leading order.

Next, one also needs to include the nonlinear terms from (2.1). This was done for DHB in the one-dimensional CGL PDE in Section 6 of [17]. There, we first used the same type of integrating factor (as used in (2.3) above) to rewrite the nonlinear PDE for $B(x, \mu)$. Then, we split the solution into two parts: $B(x, \mu) = B_p(x, \mu) + b(x, \mu)$, where we recall that B_p is the particular solution of the linearized equation. The PDE for the remainder term $b(x, \mu)$ was converted into an integral equation. We showed formally that there is a mild solution of that integral equation, using an iterative method, and that the magnitude of $b(x, \mu)$ remains small at least until $\mu = \omega_0$. Hence, the solution of the full nonlinear PDE remains near the repelling QSS at least until ω_0 .

As the estimates of the mild formulation of the one-dimensional nonlinear equation only used $L^\infty \rightarrow L^\infty$ estimates for the heat semi-group, we expect that a similar formal

analysis will hold for two spatial dimensions, as well. That is, based on decomposing $B(x, y, \mu) = B_p(x, y, \mu) + b(x, y, \mu)$, we expect that $b(x, y, \mu)$ remains small at least until μ reaches ω_0 . Fundamentally, the nonlinear terms remain exponentially small at least as long as the linear terms do.

3. The spatio-temporal memory surface and spatio-temporal buffer surface for the CGL equation

In this section, we derive the general formulas for the spatio-temporal memory and buffer surfaces of (2.1), and we apply these to several classes of initial data $A_0(x, y)$ and several different source terms $I_a(x, y)$, respectively.

3.1. The spatio-temporal memory surface By writing the homogeneous solution $A_h(x, y, \mu)$ as a single exponential function, we define the spatio-temporal memory surface to be the set of points $(x, y, \mu_{\text{mem}}(x, y))$ at which the real part of the argument of the exponential vanishes. In this section, we examine three different types of initial data: constant, Gaussian and periodic, in order to study how the spatio-temporal memory surface depends on the functional form of $A_0(x, y)$.

For constant initial data $A_0(x, y) = 1$, one finds from (2.4) that

$$A_h(x, y, \mu) = e^{(\mu^2 - \mu_0^2 + 2i\omega_0(\mu - \mu_0))/2\varepsilon}. \quad (3.1)$$

It is independent of (x, y) . Hence, $A_h(x, y)$ is exponentially small for all $\mu \in (\mu_0, -\mu_0)$. At $\mu = -\mu_0$, the real part of the exponential vanishes, which implies that the memory surface is a horizontal plane in the (x, y, μ) space: that is,

$$\mu_{\text{mem}}(x, y) = -\mu_0 \quad \text{for all } (x, y), \quad (3.2)$$

where we recall that $\mu_0 < 0$. Then, for $\mu > -\mu_0$, A_h becomes exponentially large.

For Gaussian initial data $A_0(x, y) = e^{-(x^2 + y^2)/4\sigma}$, formula (2.4) shows that the homogeneous solution is given by

$$A_h(x, y, \mu) = \left[\frac{\sigma}{\sigma + d(\mu - \mu_0)} \right] e^{(\mu^2 - \mu_0^2 + 2i\omega_0(\mu - \mu_0))/2\varepsilon} e^{-(x^2 + y^2)/(4(\sigma + d(\mu - \mu_0)))}. \quad (3.3)$$

At each (x, y) , the real part of the argument of the exponential vanishes for

$$\begin{aligned} \frac{\mu^2 - \mu_0^2}{2\varepsilon} + \ln(\sigma) - \frac{1}{2} \ln((\sigma + d_R(\mu - \mu_0))^2 + d_I^2(\mu - \mu_0)^2) \\ - \frac{(x^2 + y^2)(\sigma + d_R(\mu - \mu_0))}{4[(\sigma + d_R(\mu - \mu_0))^2 + d_I^2(\mu - \mu_0)^2]} = 0, \end{aligned} \quad (3.4)$$

where we recall that $d = d_R + id_I$ in (2.1). Hence, to leading order asymptotically, we find that the spatio-temporal memory surface is parabolic in x and y ,

$$\begin{aligned}\mu_{\text{mem}}(x, y) = |\mu_0| - \frac{\varepsilon}{|\mu_0|} \left(\ln(\sigma) - \frac{1}{2} \ln((\sigma + 2d_R|\mu_0|)^2 + 4d_I^2|\mu_0|^2) \right) \\ + \frac{\varepsilon}{4|\mu_0|} \left(\frac{(x^2 + y^2)(\sigma + 2d_R|\mu_0|)}{(\sigma + 2d_R|\mu_0|)^2 + 4d_I^2|\mu_0|^2} \right) + O(\varepsilon^2).\end{aligned}\quad (3.5)$$

For periodic initial data $A_0(x, y) = \cos(\pi(x - y)/L) \cos(\pi(x + y)/L)$, the homogeneous solution for $\mu > \mu_0$ is

$$A_h(x, y, \mu) = e^{(\mu^2 - \mu_0^2 + 2i\omega_0(\mu - \mu_0))/2\varepsilon} e^{-4\pi^2 d(\mu - \mu_0)/L^2} \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y). \quad (3.6)$$

Hence, at any point (x, y) , the real part of the argument vanishes for

$$\frac{1}{2\varepsilon}(\mu^2 - \mu_0^2) - \frac{4\pi^2 d(\mu - \mu_0)}{L^2} + \ln \left| \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y) \right| = 0. \quad (3.7)$$

Asymptotically to leading order, one finds logarithmic dependence on A_0 ,

$$\mu_{\text{mem}}(x, y) = |\mu_0| + \frac{8\varepsilon\pi^2 d}{L^2} - \frac{\varepsilon}{|\mu_0|} \ln \left| \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y) \right|. \quad (3.8)$$

These theoretical results for $\mu_{\text{mem}}(x, y)$ with constant, Gaussian and periodic initial data are compared with simulations of (2.1) in Section 4.

REMARK 1. In [17], for the CGL PDE in one dimension, we used the label homogeneous exit time curve. (There, homogeneous referred to the curve being defined by A_h , not that the curve is spatially homogeneous.) Here, we label it instead the spatio-temporal memory surface, since it is determined by the memory of the initial data. One could also label it the spatio-temporal way-in way-out surface, in analogy with the way-in way-out function defined for DHB in analytic ODEs [35, 36].

3.2. The spatio-temporal buffer surface In this section, we focus on the particular solution $A_p(x, y, \mu)$ with special emphasis on the fourth (last) term in the general formula (2.5). We consider $\delta \leq \mu \leq \omega_0$ and label this term G , where

$$G(x, y, \mu) = (\sqrt{2\pi}g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}))e^{(\mu + i\omega_0)^2/2\varepsilon}. \quad (3.9)$$

The function G is the component of the particular solution $A_p(x, y, \mu)$ that measures the deviation from the repelling QSS (where we recall that the first three terms in (2.5) represent the QSS). It is generated by passage over the saddle point at $\mu = -i\omega_0$ in the complex μ -plane, as shown in Appendix A. It is present in all solutions that start from initial data at any $\mu_0 < 0$, including solutions on the attracting QSS.

The spatio-temporal buffer surface is defined as the surface along which

$$|G(x, y, \mu)| = 1, \quad (3.10)$$

to leading order. Here, we derive a general formula for the spatio-temporal buffer surface for smooth, bounded sources $I_a(x, y)$, and asymptotic formulas for it with constant, Gaussian, periodic and stripe sources. We see that, in general, G ceases to be

exponentially small and becomes $O(1)$ in a spatially dependent manner. This implies that all solutions with initial data given at $\mu_0 < 0$, including those on the attracting QSS, must leave an $O(1)$ neighbourhood of the repelling QSS when μ reaches the buffer surface, irrespective of how large $|\mu_0|$ is; that is, irrespective of how far in the distant past the solutions approached the attracting QSS.

From (3.9), we find that the spatio-temporal buffer surface is given implicitly by

$$\operatorname{Re}\left[\ln(\sqrt{2\pi}g(x, y, \mu + i\omega_0)) + \frac{1}{2\varepsilon}(\mu + i\omega_0)^2\right] = 0$$

for general smooth and bounded source terms $I_a(x, y)$. We label the solution $\mu_{\text{buf}}(x, y)$. Its graph is the spatio-temporal buffer surface. To leading order,

$$\mu_{\text{buf}}^2(x, y) = \omega_0^2 - \varepsilon \ln(2\pi) - 2\varepsilon \ln|g(x, y, \omega_0 + i\omega_0)|. \quad (3.11)$$

At each point (x, y) , the buffer surface $\mu_{\text{buf}}(x, y)$ determines the time at which the deviation of $A_p(x, y, \mu)$ from the repelling QSS ceases to be exponentially small.

We begin with constant source terms, $I_C(x, y) = c$, and set $c = 1$ without loss of generality. Substituting this into the integral for g in (2.5), we find

$$g(x, y, \mu - \tilde{\mu}) = 1 \quad \text{for all } (x, y) \quad \text{and} \quad \mu > \tilde{\mu} \geq \mu_0. \quad (3.12)$$

Hence, by (3.11), the buffer surface is to leading order

$$\mu_{\text{buf}}(x, y) = \omega_0 - \frac{\varepsilon}{2\omega_0} \ln(2\pi) \quad \text{for all } (x, y). \quad (3.13)$$

It is independent of (x, y) . Hence, at all points in the domain, the particular solution ceases to be exponentially small uniformly at this time, μ_{buf} .

Next, we consider Gaussian source terms, $I_G(x, y) = e^{-(x^2+y^2)/4\sigma}$ with $\sigma > 0$. Using (2.5), we find

$$g(x, y, \mu - \tilde{\mu}) = \left(\frac{\sigma}{\sigma + d(\mu - \tilde{\mu})}\right) \exp\left[\frac{-(x^2 + y^2)}{4(\sigma + d(\mu - \tilde{\mu}))}\right].$$

Hence, (3.11) implies that, to leading order, the buffer surface is

$$\begin{aligned} \mu_{\text{buf}}(x, y) = \omega_0 - \frac{\varepsilon}{2\omega_0} \ln\left(\frac{2\pi\sigma^2}{(\sigma + \omega_0(d_R - d_I))^2 + \omega_0^2(d_R + d_I)^2}\right) \\ + \frac{\varepsilon}{4\omega_0} \left[\frac{(\sigma + \omega_0(d_R - d_I))}{(\sigma + \omega_0(d_R - d_I))^2 + \omega_0^2(d_R + d_I)^2}\right] (x^2 + y^2). \end{aligned} \quad (3.14)$$

For periodic source terms, $I_P(x, y) = \cos(\pi(x - y)/L) \cos(\pi(x + y)/L)$, (2.5) yields

$$g(x, y, \mu - \tilde{\mu}) = e^{-4\pi^2 d(\mu - \tilde{\mu})/L^2} \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y).$$

Hence, (3.11) implies that, to leading order, the buffer surface is

$$\begin{aligned} \mu_{\text{buf}}(x, y) = \omega_0 - \frac{\varepsilon}{2\omega_0} \ln(2\pi) + \frac{4\varepsilon\pi^2(d_R - d_I)}{L^2} \\ - \frac{\varepsilon}{\omega_0} \ln \left| \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y) \right|. \end{aligned} \quad (3.15)$$

Finally, we let H denote the Heaviside step function (1 on $x > 0$ and 0 on $x < 0$) and study stripe source terms

$$I_S(x, y) = \sum_k [H(x - x_k) - H(x - (x_k + h))].$$

Here, h is the width of the stripe, x_k denotes the left edge of the k th stripe and $x_{k+1} - x_k = \Delta > h$ for each k . We use (2.5) to derive

$$g(x, y, \mu - \tilde{\mu}) = \frac{1}{2} \sum_k \left(\operatorname{erf} \left(\frac{x_k + h - x}{\sqrt{4d(\mu - \tilde{\mu})}} \right) - \operatorname{erf} \left(\frac{x_k - x}{\sqrt{4d(\mu - \tilde{\mu})}} \right) \right), \quad (3.16)$$

where $\operatorname{erf} z = 2/\sqrt{\pi} \int_0^z e^{-t^2} dt$ is the error function. Hence, (3.11) implies that, to leading order, the buffer surface is given by

$$\mu_{\text{buf}}(x, y) = \omega_0 - \frac{\varepsilon}{2\omega_0} \ln(2\pi) - \frac{\varepsilon}{\omega_0} \ln |g(x, y, \omega_0(1 + i))|, \quad (3.17)$$

where g is given by (3.16). For these source terms, the spatio-temporal buffer surface is compared with the numerical solutions of the PDE (2.1) in Section 4.

4. DHB in the CGL equation with constant, Gaussian and stripe sources

In this section, we report on the spatially dependent duration of the DHB observed in direct numerical simulations of solutions of the CGL PDE (2.1) with $\mu_0 < 0$. We examine both cases in which the initial time satisfies $\mu_0 \in (-\omega_0, 0)$ and $\mu_0 \leq -\omega_0$. We recall that the escape surface is the graph of the time $\mu_{\text{esc}}(x, y)$ in (x, y, μ) space. Numerically, we obtain $\mu_{\text{esc}}(x, y)$ by calculating the set of points at which $|\operatorname{Re}(A_{\text{num}}(x, y)) - \operatorname{Re}(A_{\text{QSS}}(x, y))| = \delta_{\text{th}}$, where $A_{\text{num}}(x, y)$ is the numerically computed solution of (2.1), $A_{\text{QSS}}(x, y)$ is the value of A along the QSS (2.2) and δ_{th} denotes a threshold.

We show that $\mu_{\text{esc}}(x, y)$ agrees with the predictions made from the memory and buffer surfaces. At each point (x, y) ,

$$\mu_{\text{esc}}(x, y) \approx \min(\mu_{\text{mem}}(x, y), \mu_{\text{buf}}(x, y)).$$

We work with several different types of source terms (constant, Gaussian and stripe) and with the asymptotic expansions for these surfaces derived in Section 3.

In the numerical simulations, we used symmetric Strang splitting [33], with centred finite differences for the spatial discretization and fourth-order Runge–Kutta with fixed

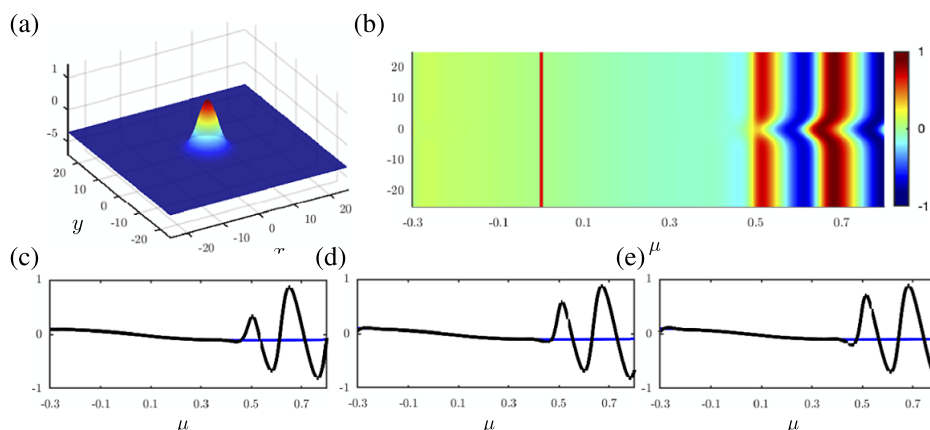


FIGURE 1. Delayed onset of oscillations in (2.1) with constant source term $I_C(x, y) = 1$ and Gaussian initial data (4.1) given at $\mu_0 = -0.3$. (a) Snapshot of $\text{Re } u$ at $\mu = -\mu_0 = 0.3$ showing that the solution is radially symmetric. (b) Space-time evolution along $y = 0$. The red line indicates the instantaneous Hopf bifurcation. The temporal evolution of $\text{Re } u$ (black curve) is compared with the QSS (blue curve) at the centre of the Gaussian at $(x, y) = (0, 0)$ (c), at a radial distance of three spatial units from the origin at $(x, y) = (3/\sqrt{2}, 3/\sqrt{2})$ (d), and at a radial distance of 10 spatial units from the origin at $(x, y) = (10/\sqrt{2}, 10/\sqrt{2})$ (e). In each case, the numerical solution stays close to the QSS well past the instantaneous Hopf bifurcation at $\mu = 0$, after which there is a hard onset to large-amplitude oscillations.

time step for the time discretization. The results were also checked independently using a Chebyshev grid for the spatial discretization, finding good agreement.

4.1. DHB with constant source term In the first representative simulation, we study DHB in (2.1) with constant source term $I_C(x, y) = 1$ and Gaussian initial data

$$A_0(x, y) = c_1 + c_2 e^{-(x^2+y^2)/4\sigma}, \quad c_1, \quad c_2 > 0, \quad (4.1)$$

given at $\mu_0 = -0.3$, which we note is in $(-\omega_0, 0)$. The dynamics are shown in Figure 1, and a quantitative comparison between the escape surface for this solution and the memory surface calculated using the analysis of the previous section is shown in Figure 2. (Note that the solution is radially symmetric, because the initial data was chosen to be radially symmetric for simplicity in this first simulation; see also Figure 1(a).)

As expected, the solution stays near the repelling QSS at least until μ reaches $-\mu_0$ (see Figure 1(a) and (b)). Then, the delayed, post-Hopf, temporal oscillations are initiated at the origin of the domain, near $\mu = -\mu_0$ (see Figure 1(c)). As μ continues to slowly increase, the oscillations occur on successively larger disks about the origin (see Figure 1(d) and (e) for two other points). Then, once μ reaches approximately 0.327 (which depends on the chosen threshold), there is a fairly rapid transition to large-scale oscillations, and these occur on the entire domain.

For this first simulation, we also show the escape surface in Figure 2(a). Below the escape surface, that is for all $O(1)$ values $\mu < \mu_{\text{esc}}(x, y)$, the solution is near the repelling QSS. Then, at each point (x, y) , the oscillations of amplitude $O(\sqrt{\varepsilon})$ set in when μ

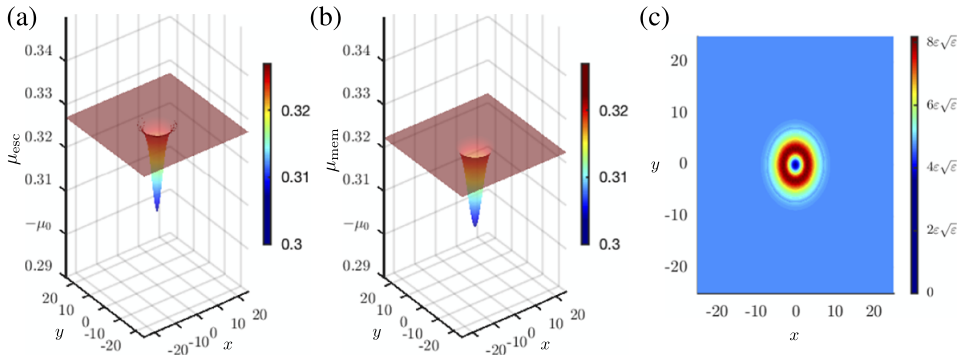


FIGURE 2. (a) In the three-dimensional (x, y, μ) space, the surface $\mu_{\text{esc}}(x, y)$ is where the hard onset of the oscillations occurs. It has been obtained from direct numerical simulation of the PDE (2.1) with constant source term $I_C(x, y) = 1$ and Gaussian initial data $A_0(x, y) = c_1 + c_2 e^{-(x^2+y^2)/4\sigma}$ with $(c_1, c_2, \sigma) = (0.5, 0.5, 2.5)$, given at $\mu_0 = -0.3$. (b) The memory surface $\mu_{\text{mem}}(x, y)$ is given by (3.4). It gives the leading order asymptotics of the escape time for all (x, y) . (c) The difference $|\mu_{\text{esc}}(x, y) - \mu_{\text{mem}}(x, y)|$ is shown in the projection onto the (x, y) plane. Here, the parameters are $d = 1$, $\omega_0 = 0.5$, $\alpha = 0.2$ and $\varepsilon = 0.01$.

reaches $\mu_{\text{esc}}(x, y)$. Furthermore, at each point (x, y) , the amplitude of the oscillations becomes large ($O(1)$) as soon as μ is slightly beyond $\mu_{\text{esc}}(x, y)$, which confirms for all points (x, y) what is shown for just three points in Figure 1.

For comparison, we show the memory surface $\mu_{\text{mem}}(x, y)$ in Figure 2(b). The memory surface was computed as follows. First, we combined the homogeneous solutions (3.1) and (3.3) to determine the homogeneous solution, A_h , with the initial data (4.1), $A_0(x, y) = c_1 + c_2 e^{-(x^2+y^2)/4\sigma}$. Then, by enforcing the condition $|A_h| = 1$, we obtained the leading order asymptotic relationship

$$\mu^2 - \mu_0^2 + 2\varepsilon \ln \left| c_1 + c_2 \frac{\sigma}{\sigma + d(\mu - \mu_0)} \exp \left(\frac{-(x^2 + y^2)}{4(\sigma + d(\mu - \mu_0))} \right) \right| = 0 \quad (4.2)$$

for the memory surface corresponding to this A_0 .

The difference $|\mu_{\text{esc}}(x, y) - \mu_{\text{mem}}(x, y)|$ is shown in Figure 2(c). At the centre of the Gaussian, the difference is small (with magnitude of approximately 10^{-5} , that is, $O(\varepsilon^{5/2})$). Then, in an annular region about the origin (red and dark red), the difference is slightly larger, due to nonlinear effects. Finally, both surfaces exhibit a fairly rapid transition into the regime (blue region in Figure 2(c), red in (a) and (b)) where they are essentially constant, since the Gaussian is tiny. Here, $\mu_{\text{mem}}(x, y) = \sqrt{\mu_0^2 - 2\varepsilon \ln c_1} \approx 0.322$ to leading order (as obtained from (4.2) for large $x^2 + y^2$), which agrees well with the value observed numerically in (2.1). Similar results were observed for solutions and the escape and memory surfaces with other values of $-\omega_0 < \mu_0 < 0$ (data not shown).

Moreover, for the CGL with this (constant) source term, the buffer surface lies above the memory surface for all (x, y) in three dimensions, since

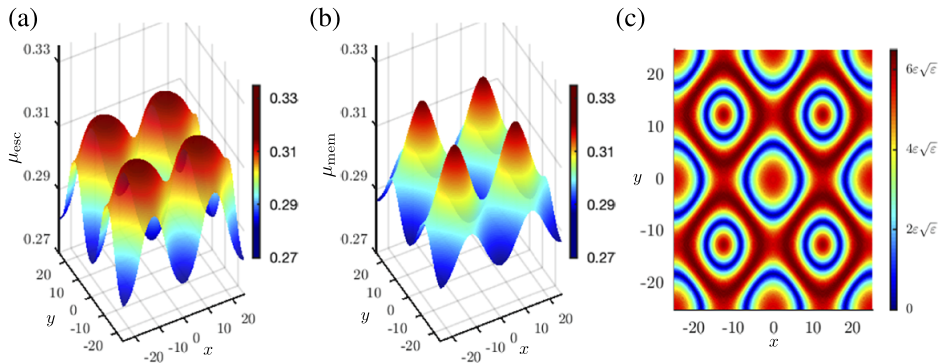


FIGURE 3. (a) In the three-dimensional (x, y, μ) space, the surface $\mu_{\text{esc}}(x, y)$ is where the hard onset of the oscillations occurs. It has been obtained from direct simulation of (2.1) with constant source term $I_C(x, y) = 1$ and periodic initial data $A_0(x, y) = p_1 + p_2 \cos(\pi(x - y)/L) \cos(\pi(x + y)/L)$ at $\mu_0 = -0.3$ with $(p_1, p_2, L) = (1, 0.5, 25)$. (b) The memory surface $\mu_{\text{mem}}(x, y)$ given by (3.7). (c) The surface $|\mu_{\text{num}}(x, y) - \mu_{\text{mem}}(x, y)|$ shown in the projection onto the (x, y) plane. The parameters are $d = 1$, $\omega_0 = 0.5$, $\alpha = 0.2$ and $\varepsilon = 0.01$.

$\mu_{\text{buf}}(x, y) = \sqrt{\omega_0^2 - \varepsilon \ln(2\pi)} \approx 0.481$ for all (x, y) . Hence, for all (x, y) , the minimum in $\mu_{\text{esc}}(x, y) \sim \min\{\mu_{\text{mem}}(x, y), \mu_{\text{buf}}(x, y)\}$ is given entirely by the memory surface in this simulation.

The second representative simulation is also with $I_C(x, y) = 1$. However, now the initial data given at $-\omega_0 < \mu_0 < 0$ is periodic,

$$A_0(x, y) = p_1 + p_2 \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y),$$

with $p_1 > p_2 > 0$, so that $A_0(x, y)$ is strictly positive everywhere. The results are shown in Figure 3.

The escape surface computed from the numerical simulations (Figure 3(a)) shows that the oscillations first occur at time $\mu \approx 0.281$ at the points $(x, y) = (j_1 L, j_2 L)$, where $j_1, j_2 = -1, 0, 1$ (dark blue), that is, at the maxima of $A_0(x, y)$, where $\cos(\pi(x - y)/L) \cos(\pi(x + y)/L) = 1$. About each of those points, the oscillations set in on successively larger disks as μ slowly increases, until those disks collide at $\mu \approx 0.306$ (near the transition from yellow to orange). Then, as μ increases further, the escape surface consists of the four inverted paraboloid segments (orange and red). The peaks occur at time $\mu \approx 0.315$ at the points $(x, y) = (j_1 L/2, j_2 L/2)$, where $j_1, j_2 = -1, 1$ (dark red), that is, at the minima of $A_0(x, y)$, where $\cos(\pi(x - y)/L) \cos(\pi(x + y)/L) = -1$.

For comparison, the memory surface $\mu_{\text{mem}}(x, y)$ is defined by

$$\mu^2 - \mu_0^2 + 2\varepsilon \ln \left| p_1 + p_2 \exp \left(-\frac{4\pi^2 d(\mu - \mu_0)}{L^2} \right) \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y) \right| = 0, \quad (4.3)$$

and shown in Figure 3(b). It has global minima (dark blue) when

$$\mu_{\text{mem}} = \sqrt{|\mu_0|^2 - 2\varepsilon \ln(1 + 0.5e^{(-8\pi^2 d|\mu_0|)/L^2})} \approx 0.2866$$

to leading order, with $\mu_0 = -0.3$. These occur at the points where $A_0(x, y)$ has its maxima. Also, the memory surface has global maxima (dark red) when $\mu \approx 0.3211$, at the points where $A_0(x, y)$ has its minima. In between, it has the same conical shape qualitatively as the escape surface.

The difference between $\mu_{\text{esc}}(x, y)$ and $\mu_{\text{mem}}(x, y)$ is shown in Figure 3(c). The nonlinear terms cause μ_{esc} to grow more steeply from the local minima, compared with the memory surface, so that the disks collide at a slightly later time than for the memory surface. Then, after the disks collide, the inverted paraboloids are slightly wider in $\mu_{\text{esc}}(x, y)$ due to the nonlinear terms.

Finally, for this second representative simulation, we report that the buffer surface also lies above the memory surface for all (x, y) in three dimensions. Indeed, with constant source $I_C(x, y) = 1$, $\mu_{\text{buf}}(x, y) = \sqrt{\omega_0^2 - \varepsilon \ln(2\pi)} \approx 0.481$. Hence, here also the minimum is given by the memory surface for all (x, y) .

REMARK 2. In the second simulation, we also combined two homogeneous solutions, (3.1) and (3.6), to determine the homogeneous solution A_h that corresponds to the periodic initial profile $A_0(x, y) = p_1 + p_2 \cos(\pi(x - y)/L) \cos(\pi(x + y)/L)$. Then, by setting $|A_h| = 1$, we obtain (4.3).

REMARK 3. We also explored the effect of giving the initial data at different times, including $\mu_0 < -\omega_0$, while keeping the source term and initial data the same as in the second simulation. In these simulations (data not shown), we observe that $\mu_{\text{esc}}(x, y) = \omega_0 - \varepsilon \ln(2\pi)$ to leading order for all (x, y) , as determined by A_p (recall (3.13)). This is as expected from the analysis, because $\mu_{\text{buf}}(x, y) < \mu_{\text{mem}}(x, y)$ at all points when the initial data is given at time $\mu_0 < -\omega_0$ and ε is sufficiently small.

4.2. DHB with Gaussian source term Gaussian source terms can be used to model spatially localized, radially symmetric inputs, such as a circular spot of visible light that shines on the reactor in a chemical pattern-forming experiment. A representative example of (2.1) with Gaussian source term is shown in Figure 4. Here, $I_G(x, y) = \exp(-(x^2 + y^2)/4\sigma)$ with $\sigma = 1$, and the initial data at $\mu_0 = -0.75$ (which is chosen so that $\mu_0 < -\omega_0$) is

$$A_0(x, y) = \cos \frac{\pi}{L}(x - y) \cos \frac{\pi}{L}(x + y) \quad \text{with } L = 25.$$

From direct numerical simulations of the PDE, we find that the domain of the escape surface (Figure 4(a)) can be split into two distinct parts: the annular region $R_G = \{(x, y) \mid x^2 + y^2 \leq 11.5^2\}$ and its complement $R_P = \mathbb{R}^2 \setminus R_G$. On R_G , the earliest numerically detected escape occurs at the origin for $\mu \approx 0.4905$. The escape surface is radially symmetric and increases on concentric rings, which thus creates a rotationally symmetric paraboloid. In contrast, for $(x, y) \in R_P$, the escape surface is no longer radially symmetric, and is instead a periodic tile pattern, reflecting the initial data.

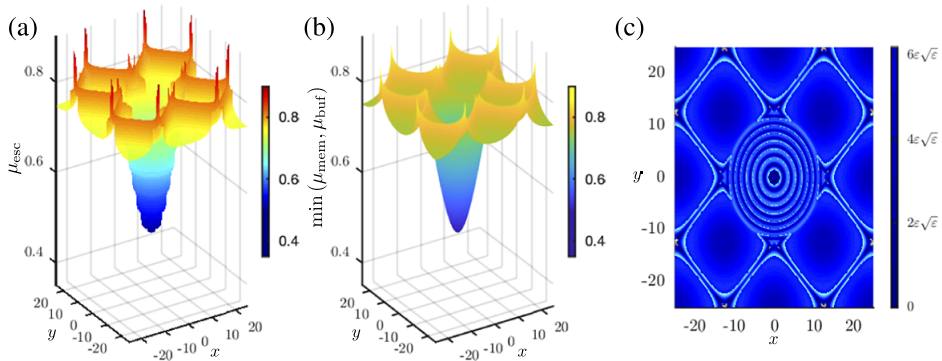


FIGURE 4. (a) In the three dimensional (x, y, μ) space, the surface $\mu_{\text{esc}}(x, y)$ is where the hard onset of the oscillations occurs. It has been obtained from direct simulation of (2.1) with Gaussian source term $I_G(x, y) = \exp(-(x^2 + y^2)/4\sigma)$ and periodic initial data $A_0(x, y) = \cos(\pi(x - y)/L) \cos(\pi(x + y)/L)$ at $\mu_0 = -0.75$ with $L = 25$. (b) The predicted escape surface is given by the minimum of $\mu_{\text{buf}}(x, y)$ (parabolic part) and $\mu_{\text{mem}}(x, y)$ (periodic part). (c) The surface $|\mu_{\text{esc}}(x, y) - \min\{\mu_{\text{mem}}, \mu_{\text{buf}}\}|$ shown in the projection onto the (x, y) plane. The parameters are $d = 1$, $\omega_0 = 0.5$, $\alpha = 0.2$ and $\varepsilon = 0.01$.

The minima occur at $\mu \approx 0.738$ (yellow/green regions) and the maxima at $\mu \approx 0.812$ (red lines between yellow/green regions).

For this source term and initial condition, the buffer surface is given by (3.14) and the memory surface is given by (3.7). In Figure 4(b), we plot $\min\{\mu_{\text{buf}}(x, y), \mu_{\text{mem}}(x, y)\}$. On R_G , the minimum is given by the buffer surface, whereas it is given by the memory surface on R_P . The buffer surface predicts that the earliest onset occurs at the origin at $\mu \approx 0.4908$. The values of μ_{buf} increase in a radially symmetric fashion until $\mu \approx -\mu_0 = 0.75$ on the boundary of R_G . Then, for (x, y) on R_P , the memory surface predicts the onset. (Compare Figure 4(a) and (b).)

The difference $|\mu_{\text{esc}}(x, y) - \min\{\mu_{\text{buf}}(x, y), \mu_{\text{mem}}(x, y)\}|$ is small throughout, as shown in Figure 4(c). The difference is especially small (of the order of the neglected terms in the asymptotic expansion) near the onset at the origin, which is at the minimum of the buffer surface, and also near the local minima of the memory surface.

4.3. DHB with stripe source term Stripe source terms are also of fundamental interest. For example, chemical reactions irradiated with constant intensity light filtered through stripe masks can produce complex patterns (see [7, 16] and references within).

Here, we use a simple model for a strictly positive stripe source term,

$$I_S(x, y) = 1 + \frac{1}{2} \sum_{k=-2}^2 [H(x - x_k) - H(x - x_k - h)], \quad (4.4)$$

where h is the stripe width, x_k denotes the left edge of the k th stripe and $x_{k+1} - x_k = \Delta > h$. We set $x_0 = -1.25$, $h = 2.5$ and $\Delta = 10$.

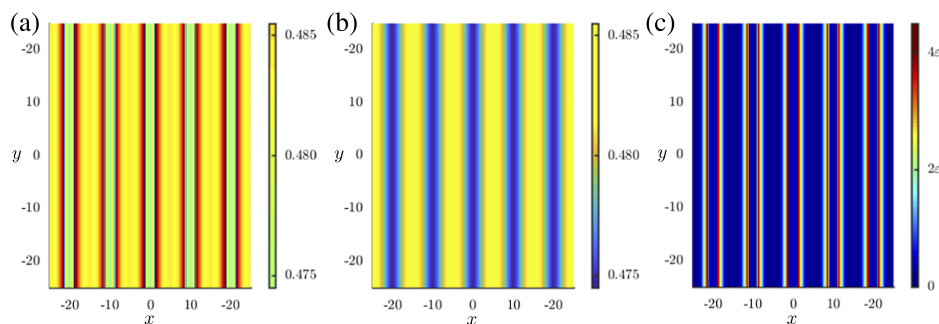


FIGURE 5. (a) The surface $\mu_{\text{esc}}(x, y)$, shown here in the projection onto the (x, y) plane, is where the hard onset of oscillations occurs. It has been obtained from numerical simulation of (2.1) with stripe source term $I_S(x, y)$ (4.4) and constant initial data $A_0(x, y) = 1$ given at $\mu_0 = -1$. (b) The buffer surface $\mu_{\text{buf}}(x, y)$ defined by (3.11), with g given by the linear combination of (3.12) and (3.16). (c) The difference $|\mu_{\text{esc}}(x, y) - \mu_{\text{buf}}(x, y)|$. The parameters are $d = 1$, $\omega_0 = 0.5$, $\alpha = 0.2$ and $\varepsilon = 0.01$.

For the PDE (2.1), the surfaces $\mu_{\text{esc}}(x, y)$ (calculated from direct numerical simulations) and $\mu_{\text{buf}}(x, y)$ (calculated from (3.12) and (3.16)) are shown in Figure 5(a) and (b), along with the difference in Figure 5(c). From the plot of $\mu_{\text{esc}}(x, y)$, we see that the hard onset of oscillations occurs first at the points where $I_S(x, y)$ has its maximum, that is, inside the stripes. See the green stripes in the plot of $\mu_{\text{esc}}(x, y)$ in Figure 5(a). The buffer curve $\mu_{\text{buf}}(x, y)$ also has local minima inside the stripes (blue stripes in Figure 5(b)), and these stripes have approximately the same width as those of the escape surface. The difference in the green and blue stripes is of $O(\varepsilon\sqrt{\varepsilon})$.

In the complementary regions, the PDE simulations show that the duration of the DHB is longer. The escape surface lies just below ω_0 , by about 2ε (see the yellow stripes in Figure 5(a)). In the complementary regions, the buffer surface transitions more gradually to its local maxima, which also occur at the local minima of the source term. Here, the difference is of $O(\varepsilon\sqrt{\varepsilon})$.

The narrow red strips between the green and yellow bands in Figure 5(a) correspond to the regions where the Heaviside function has a discontinuity. This discontinuity is smoothed out in the predicted buffer surface Figure 5(b), so that $\mu_{\text{buf}}(x, y)$ is smooth and continuous across the whole domain. This difference is highlighted in Figure 5(c). There, the error is small and of $O(\varepsilon\sqrt{\varepsilon})$ throughout the blue regions, and is of $O(\varepsilon)$ in the thin strips where the Heaviside function has jump discontinuities.

We add that, in this simulation with the stripe source term (4.4) and the constant initial data given at $\mu_0 = -1$, the memory surface (3.2) lies well above the buffer surface (3.13) for all (x, y) . Hence, the escape time is determined exclusively by the buffer surface, as reported.

4.4. The competition between $\mu_{\text{mem}}(x, y)$ and $\mu_{\text{buf}}(x, y)$ In the above simulations, we have shown that at each point (x, y) , $\mu_{\text{esc}}(x, y) \sim \min(\mu_{\text{mem}}(x, y), \mu_{\text{buf}}(x, y))$ to leading

order. Therefore, in effect, there is a competition at every point (x, y) between $A_h(x, y, \mu)$ and $A_p(x, y, \mu)$ in which the first one that ceases being exponentially small determines the maximal duration of the delay in the Hopf bifurcation to leading order.

The memory surface $\mu_{\text{mem}}(x, y)$ depends to leading order on the initial time μ_0 . Then, at $O(\varepsilon)$, it also depends on the logarithm of $|A_h(x, y; \mu)|$ (recall, for example, equations (3.2), (3.5) and (3.8)).

The space-time buffer surface $\mu_{\text{buf}}(x, y)$ (recall (3.11) in Section 3.2) is determined to leading order by ω_0 , which is the frequency at the instantaneous Hopf bifurcation, and then at $O(\varepsilon)$ by the logarithm of the Stokes term $g(x, y, \omega_0(1 + i))$ in (3.11). Indeed, as highlighted by the analysis in the complex μ plane in Appendix A, the Stokes line through the saddle point at $\mu = -i\omega_0$ crosses the negative μ -axis at $\mu = -\omega_0$ (see Figure A.1). Hence, the initial time μ_0 is to the left of that here, and the contour used to find $A_p(x, y, \mu)$ for values of $\mu > 0$ goes through that saddle. The saddle point method (applied in Appendix A) shows that the function $G(x, y, \mu)$ given by (3.9) arises owing to passage over the saddle at $-i\omega_0$. For each $\mu > 0$, $G(x, y, \mu)$ measures the deviation of the solution A_p from the repelling QSS. On $\mu > 0$, it stays exponentially small at least until μ reaches $+\omega_0$, where the other Stokes line through the saddle reaches the positive μ -axis. Then, as μ slowly increases beyond ω_0 , $G(x, y, \mu)$ ceases to be exponentially small—and grows to become $O(1)$ —in a spatially dependent manner. This is why the buffer surface is defined by (3.11), that is, when and where $|G| = 1$.

Therefore, at $O(1)$, μ_0 and ω_0 determine the competition, and at $O(\varepsilon)$ it is determined by the logarithmic terms. Moreover, for any finite value of ε , there can be a changeover between the two terms for which one wins the competition, since, in general, the logarithmic terms in both expressions can grow due to their spatial dependence (see, for example, Figure 4).

5. DHB in the two-dimensional Brusselator model

In two space dimensions, the Brusselator model with a source term and a slowly varying rate constant is given by the system

$$\begin{aligned} u_t &= a(x, y) - (1 + b)u + u^2v + \varepsilon d_u \Delta u, \\ v_t &= bu - u^2v + \varepsilon d_v \Delta v, \\ b_t &= \varepsilon. \end{aligned} \tag{5.1}$$

Here, the independent variables are $(x, y) \in \mathbf{R}^2$ and $t \geq 0$, $\Delta = (\partial^2/\partial x^2) + (\partial^2/\partial y^2)$, and the subscript t denotes the partial derivative on t . The dependent variable u denotes the concentration of the activator chemical and v denotes that of the inhibitor species. The small parameter $0 < \varepsilon \ll 1$ measures the separation in the time scales. The spatially dependent source term, $a(x, y)$, is taken to be positive, bounded, $O(1)$ with respect to ε and smooth. The parameter $b > 0$ denotes a rate constant that slowly increases in time, and $d_u, d_v > 0$ are the diffusivities of the activator and inhibitor.

5.1. Hopf bifurcation of the QSS The Brusselator PDE (5.1) has a QSS,

$$\begin{aligned} u_{\text{QSS}}(x, y) &= a(x, y) + \varepsilon \Delta \left(d_u a(x, y) + d_v \frac{b}{a(x, y)} \right) + O(\varepsilon^2), \\ v_{\text{QSS}}(x, y) &= \frac{b}{a(x, y)} + \frac{\varepsilon}{a^2(x, y)} \Delta \left((1 - b) \frac{d_v b}{a(x, y)} - b d_u a(x, y) \right) + O(\varepsilon^2). \end{aligned}$$

We perform a linear stability analysis about the QSS by setting $u = u_{\text{QSS}} + \hat{u}$ and $v = v_{\text{QSS}} + \hat{v}$. The linearized system for \hat{u} and \hat{v} is

$$\begin{bmatrix} \hat{u}_t \\ \hat{v}_t \end{bmatrix} = M \begin{bmatrix} \hat{u} \\ \hat{v} \end{bmatrix} + \varepsilon \begin{bmatrix} d_u & 0 \\ 0 & d_v \end{bmatrix} \begin{bmatrix} \Delta \hat{u} \\ \Delta \hat{v} \end{bmatrix},$$

where

$$M = \begin{bmatrix} 2u_{\text{QSS}}v_{\text{QSS}} - (1 + b) & u_{\text{QSS}}^2 \\ b - 2u_{\text{QSS}}v_{\text{QSS}} & -u_{\text{QSS}}^2 \end{bmatrix}.$$

Hence, by setting $\text{Tr}(M) = 0$, we find that the PDE (5.1) exhibits a spatially dependent Hopf bifurcation at $b = b_H(x, y)$, where

$$b_H(x, y) = \frac{1 + a^2 + 2\varepsilon a d_u \Delta a + O(\varepsilon^2)}{1 + 2\varepsilon d_v((1 - a^2)/a)\Delta(1/a) + O(\varepsilon^2)}, \quad (5.2)$$

and we have written a for $a(x, y)$. At each point $(x, y) \in \mathbf{R}^2$, the QSS is linearly stable for $b < b_H(x, y)$ and linearly unstable for $b > b_H(x, y)$. Moreover, the frequency, ω_H , at b_H is $\omega_H = a(x, y) + O(\varepsilon)$.

REMARK 4. The Brusselator model (5.1) exhibits (super-critical) Turing bifurcations. For example, with source $a(x, y) = a_0 + \varepsilon a_1(x, y) + O(\varepsilon^2)$, where $a_0 > 0$ and $O(1)$, Turing bifurcations occur at $b_T = (1 \pm \sqrt{d_u/d_v} a_0)^2 + O(\varepsilon)$, in which spatially periodic states are created. The parameter values here were chosen so the Turing bifurcations do not have an impact on the DHB phenomenon. It would be of interest to study the interactions between the two bifurcations.

5.2. Spatio-temporal dependence of the DHB in the Brusselator model In this section, we present the results of numerical simulations showing the spatio-temporal dependence of the DHB in the Brusselator model (5.1). We take the initial values $b(t_0) < b_H(x, y)$ for all (x, y) , so that the system slowly passes through the Hopf bifurcation $b_H(x, y)$ as b slowly increases (recall (5.2)). We work with a stripe source term,

$$a_S(x, y) = 1 + \varepsilon \alpha \sum_k [H(x - x_k) - H(x - (x_k + h))], \quad (5.3)$$

where $\alpha > 0$, x_k denotes the left edge of the k th stripe, $x_{k+1} - x_k = \Delta$, $h < \Delta$ is the width of the stripe and the stripes are sufficiently well separated.

In the simulations, the solutions rapidly approach the attracting QSS for $b < b_H(x, y)$ and, after the instantaneous Hopf point is crossed, they stay near the repelling QSS for long times, of $O(1/\varepsilon)$ in the fast time t . This is the delay in the Hopf bifurcation, and the solutions leave a neighbourhood of the QSS in a spatially dependent manner. Moreover, they do so by making a large rapid jump, so that there is a hard onset of large-amplitude oscillations.

For the representative simulation shown in Figure 6, we set $\alpha = 5$, $h = 0.1$, $x_0 = -0.05$ and $\Delta = 0.3$. The initial data is spatially periodic,

$$u_0(x, y) = a(x, y) + 0.5 \cos(\pi^2 xy), \quad v_0(x, t) = \frac{b_0}{a(x, y)} + \varepsilon. \quad (5.4)$$

The initial value of the parameter is $b_0 = b(t_0) = 1.5$ (Figure 6(a)), below the Hopf point $b_H = 2$ here. As b slowly increases, the solution rapidly approaches the attracting QSS, and the component of the solution given by the spatially periodic initial data decreases. By the time $b = 1.7$, the remnant of the initial data is faint (Figure 6(b)). Then, when b reaches $b_H = 2$, there is no visible trace of the initial data (Figure 6(c)).

As b continues to increase, the solution remains near the repelling QSS (Figure 6(d), where $b = 2.2$). It is not until $b = 2.4$ that small amplitude temporal (post-Hopf) oscillations set in (Figure 6(e)). Also, the solution has regained a significant component given by the initial data. The oscillations rapidly gain large amplitude as b increases further (Figure 6(f)–(h), where $b = 2.6$, $b = 2.8$, $b = 3.0$; note the changes in the vertical scales). Hence, the observed onset is a hard onset. Also, significant interactions occur between the components coming from the initial data and the stripe source.

Further increases in b result in spatio-temporal patterns, such as spiral waves (Figure 6(i)). These arise from the interaction of the oscillations (which occur beyond the memory surface), the stripe source and the diffusion.

A cross-section through $y = 0$ is shown in Figure 6(j). The space-time plot shows that the solution stays close to the QSS (blue state) well past the instantaneous Hopf bifurcation (red line) and the large-amplitude oscillations do not set in until $b \approx 2.5$. Cross-sections taken at other values of y are qualitatively the same, with some slight shifts in x .

The hard onset of oscillations occurs along the escape surface, $b_{\text{esc}}(x, y)$, which at each point (x, y) is given by the value of b at which

$$|(u(x, y), v(x, y)) - (u_{\text{QSS}}(x, y), v_{\text{QSS}}(x, y))| = \delta_{\text{th}}.$$

This is the analogue of $\mu_{\text{esc}}(x, y)$. Plots of $b_{\text{esc}}(x, y)$ are shown for two different initial values: $b(t_0) = 0.1$ in Figure 7(a) and $b(t_0) = 1.5$ in Figure 7(b). Here, the parameters are the same as in Figure 6. These escape surfaces reveal that the spatio-temporal buffer surface (defined by the particular solution of the equation linearized about the QSS) and the spatio-temporal memory surface (defined by the homogeneous solution) play similar roles here for (5.1). For initial values $b(t_0)$ far from $b_H = 2$, as in Figure 7(a), the escape surface is determined entirely by the stripe source and,

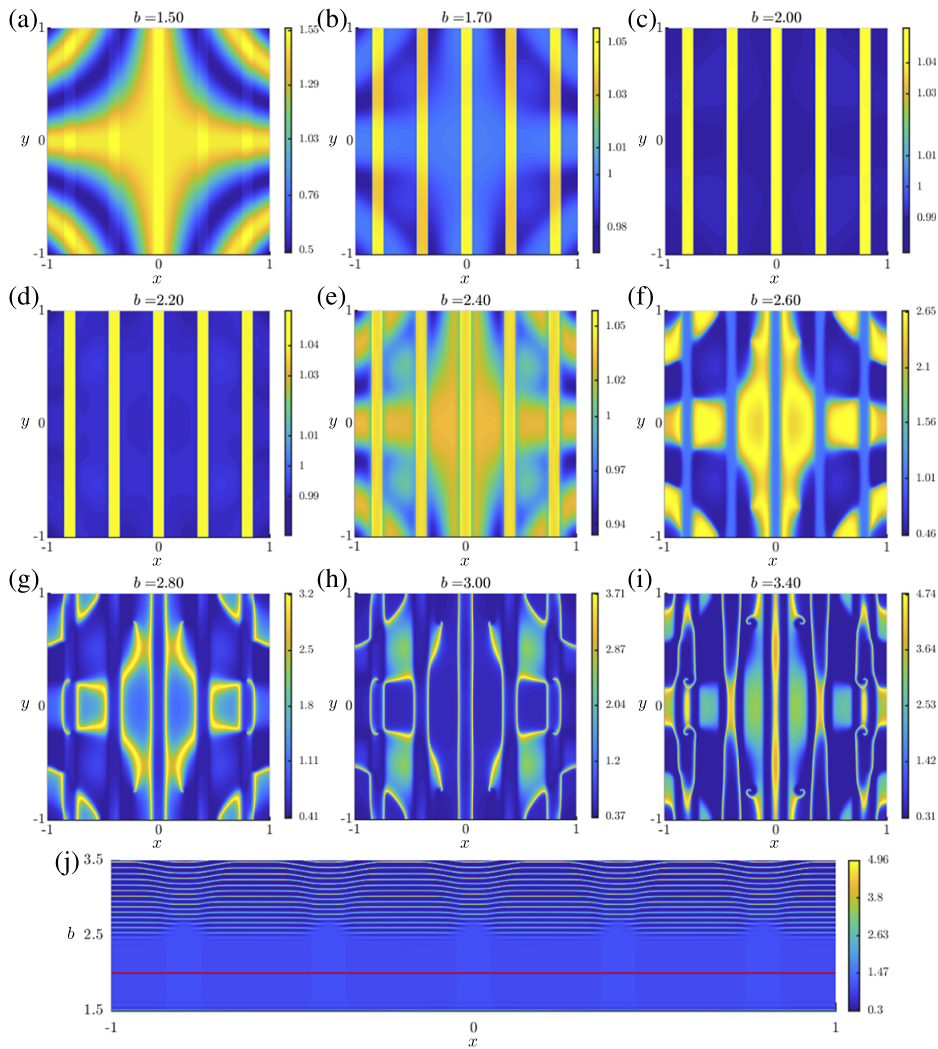


FIGURE 6. DHB in the two-dimensional Brusselator with stripe source (5.3) and spatially periodic initial data (5.4) given at $b(t_0) = 1.5$. (a)–(i) The solution $u(x, y)$ rapidly converges to the source-dependent QSS, stays close to the QSS well beyond the instantaneous Hopf bifurcation at $b = 2$ and exhibits a hard transition to oscillations. Note the different scales on the vertical colour bars in panels (a)–(i). (j) Space-time evolution along $y = 0$. The parameters are $\varepsilon = 0.01$, $d_u = 1 \times 10^{-3}$, $d_v = 5 \times 10^{-4}$, $\alpha = 5$, $h = 0.1$, $x_0 = -0.05$ and $\Delta = x_{k+1} - x_k = 0.4$.

hence, by the particular solution. The initial data has no influence on $b_{\text{esc}}(x, y)$, because the escape surface is almost invariant. Indeed, this was verified across four different choices of $b(t_0)$. Then, for initial values of $b(t_0)$ close to b_H (as in (b)), the escape surface is located below the buffer surface at all points. The homogeneous solution

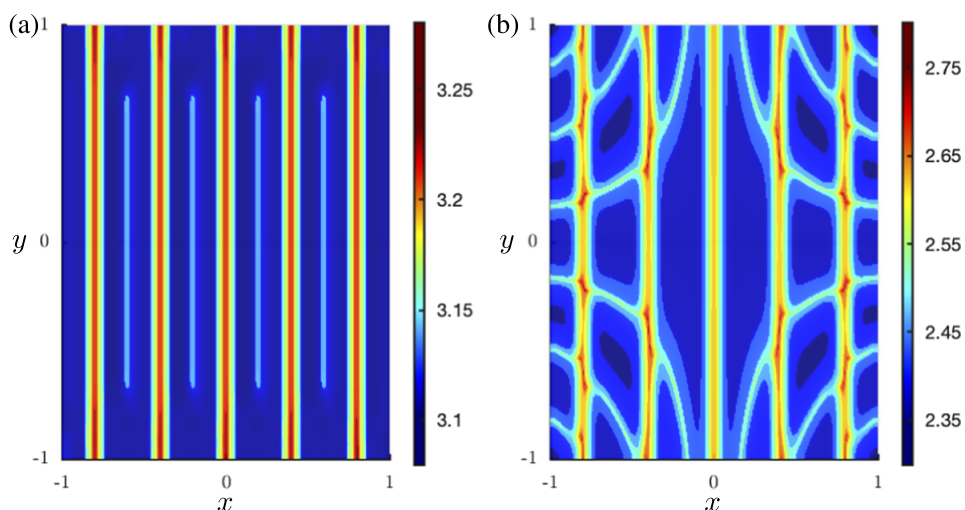


FIGURE 7. Contours of the escape surface, $b_{\text{esc}}(x, y)$, where the hard onset of oscillations occurs. The parameters are as in Figure 6 with different values of b at the initial time t_0 : $b(t_0) = 0.1$ (a) and $b(t_0) = 1.5$ (b). The escape surface in (a) depends only on the stripe source and has no memory of the initial data. Similar escape surfaces to (a) are observed for $b(t_0) = 0, 0.2$ and 0.3 . The escape surface in (b) depends on the initial profile.

determines the exit time for all (x, y) . The earliest escape times (dark blue regions) are where the initial data has its smallest amplitude (in the hyperbolic bands with the minima of the cosine term) and the light blue hyperbolic bands occur where the cosine term is maximal. In the regions of the stripes, the initial data is larger and the hard onset occurs later, although still before the buffer surface.

Overall, we find that these primary features of DHBs in the two-dimensional Brusselator—the long, spatially dependent DHB; the hard, spatially dependent onset of large-amplitude oscillations; and the properties of the escape surface—are the same as those observed for DHBs in the CGL PDE in two dimensions (Section 4), as well as for DHBs in PDEs in one dimension [3, 17, 25], in analytic ODEs [4, 20, 35–39, 44, 46, 47] and in the applications cited in the Introduction.

6. Conclusions

In this paper, we studied a pair of prototypical pattern-forming PDEs in two space dimensions. In both, an attracting QSS becomes a repelling QSS as a key parameter varies slowly in time through a generic, super-critical Hopf bifurcation. We reported the discovery of DHBs in which solutions of these PDEs in two dimensions stay near the (post-Hopf) repelling QSS for long, spatially dependent periods of time ($O(1/\varepsilon)$ fast time), before the spatially dependent hard onset of large-amplitude oscillations occurs.

First, we generalized the asymptotic and numerical results of [17] for DHBs in the CGL PDE in one space dimension to two dimensions. We introduced the spatio-temporal memory surface $\mu_{\text{mem}}(x, y)$ and the spatio-temporal buffer surface $\mu_{\text{buf}}(x, y)$, and we derived asymptotic formulas for them in the CGL PDE (recall Sections 3.1 and 3.2). These are the two-dimensional analogues of the one-dimensional memory and buffer curves introduced in [17].

The memory surface $\mu_{\text{mem}}(x, y)$ is determined by when the homogeneous solution of the linearized equation reaches amplitude one, and hence stops being exponentially small. It is labelled the memory surface for the following reason. Although the memory of the initial data fades quickly after the initial time $\mu_0 < 0$, because the solution rapidly (exponentially in time) approaches the attracting QSS before the instantaneous Hopf bifurcation, the memory of it is not lost. On the contrary, the component of the solution given by the initial data re-emerges (or resurges) in a spatially dependent manner at $\mu_{\text{mem}}(x, y)$ to leading order. The asymptotic formula for $\mu_{\text{mem}}(x, y)$ quantifies its dependence on μ_0 and on the spatial structure of the general bounded, smooth initial data $A_0(x, y)$. We applied it to several structurally different types of data (see (3.2), (3.5) and (3.8)).

The buffer surface $\mu_{\text{buf}}(x, y)$ is determined from the particular solution $A_p(x, y, \mu)$ by when $|G| = 1$ (recall (3.10)). It plays a central role in the dynamics of solutions for which initial data $A_0(x, y)$ is given at $\mu_0 \leq -\omega_0$. In particular, to leading order, all solutions with initial data given at $\mu_0 \leq -\omega_0$, including those on the attracting QSS, leave an $O(1)$ neighbourhood of the repelling QSS when μ reaches the buffer surface, irrespective of how large $|\mu_0|$ is; that is, irrespective of far in the distant past the solutions approached the attracting QSS. Furthermore, we derived the asymptotic formula for $\mu_{\text{buf}}(x, y)$, finding its dependence on the frequency ω_0 at the Hopf point and on the spatial structure of general, smooth, bounded source terms. Also, we applied the general formula to the cases of constant, Gaussian, periodic and stripe sources (see (3.13), (3.14), (3.15) and (3.17)).

Overall, for solutions of the CGL PDE (2.1) with initial data given at $\mu_0 < 0$, we found that the spatially dependent duration of the DHB is given by $\mu_{\text{esc}}(x, y) = \min\{\mu_{\text{mem}}(x, y), \mu_{\text{buf}}(x, y)\}$ to leading order. Also, we found that there is good quantitative agreement between the asymptotic formulas for $\mu_{\text{mem}}(x, y)$ and $\mu_{\text{buf}}(x, y)$ and the numerically calculated escape times $\mu_{\text{esc}}(x, y)$ for different types of initial data and source terms (see Section 4).

We distinguished between the DHBs observed for solutions with initial data given at $\mu_0 \in (-\omega_0, 0)$ and for those with initial data given at $\mu_0 \leq -\omega_0$. For the former ($\mu_0 \in (-\omega_0, 0)$), there are points at which the memory surface is reached before the buffer surface. We have shown examples where the entire memory surface lies below the buffer surface in three dimensions: that is, $\mu_{\text{mem}}(x, y) < \mu_{\text{buf}}(x, y)$ for all (x, y) (recall the first and second representative simulations in Section 4.1). There are also examples in which only a portion of the memory surface lies below the buffer surface. In this case, $A_0(x, y)$ is such that $\mu_{\text{mem}}(x, y)$ grows with $\|(x, y)\|$ and then exceeds $\mu_{\text{buf}}(x, y)$ on some part of the domain, since ε has a small but finite value.

For the latter (where $\mu_0 \leq -\omega_0$), the buffer surface is reached first as μ slowly increases. That is, there are points (x, y) at which $\mu_{\text{buf}}(x, y) \leq \mu_{\text{mem}}(x, y)$ (noting that equality can arise when $\mu_0 = -\omega_0$). Crucially, there are two cases: (1) $\mu_{\text{buf}}(x, y) \leq \mu_{\text{mem}}(x, y)$ for all (x, y) and (2) $\mu_{\text{buf}}(x, y) \leq \mu_{\text{mem}}(x, y)$ in some regions while $\mu_{\text{mem}}(x, y) < \mu_{\text{buf}}(x, y)$ in the complementary regions. The example in Section 4.3 illustrates case (1), and an example of case (2) is given in Section 4.2. In case (2), the source term is such that it grows with $\|(x, y)\|$ and then exceeds $\mu_{\text{mem}}(x, y)$ on some part of the domain. In all simulations, we found $\mu_{\text{esc}}(x, y) \approx \min\{\mu_{\text{mem}}(x, y), \mu_{\text{buf}}(x, y)\}$.

Finally, we showed numerically that DHB occurs in the Brusselator model in two space dimensions. For several source terms $a(x, y)$, the solutions spend spatially dependent, $O(1/\varepsilon)$ long times near the repelling QSS. Also, the spatially dependent onset of oscillations is a hard onset, as shown, for example, in going from Figure 6(e) to (f), where we note the change in scales. Moreover, after the oscillations have set in, there are dynamically complex interactions between the source term and the re-emergent component given by the initial data.

Looking ahead, we observe that it may be possible to extend the results for the CGL PDE in two dimensions to include some source terms $I_a(x, y)$ that are not strictly positive or bounded. For example, in Section 8 of [17], we showed that the asymptotics for the buffer curve in the one-dimensional CGL PDE extend to include bounded sign-changing source terms (such as $\cos(x)$) and even some algebraically growing source terms (quadratic in space), and that the asymptotics still agree well with the results of numerical simulations in one dimension. Also, it may be possible to extend the results to include large-amplitude source terms and $O(1)$ diffusivity, as was done for the one-dimensional CGL PDE in [17, Sections 9 and 10].

We think it would be of interest to put the Brusselator PDE (5.1) into normal form for dynamic super-critical Hopf bifurcation, and to carry out an analysis of the linearized equation in the complex time plane similar to that done for the CGL PDE (2.1) in two dimensions. It would be of interest to determine where the saddle points are and where the Stokes lines through them cross the real time axis, as well as to derive the homogeneous and particular solutions of the linearized equation and hence determine the spatio-temporal memory and buffer surfaces.

The spatially dependent source terms studied here are similar to those studied in reaction–diffusion equations in two (or more) spatial variables. In [27], the authors established the conditions under which radially symmetric inhomogeneities create coherent structures, including sources, contact defects and sinks. Also, spatial inhomogeneities have been analysed in the Swift–Hohenberg and Ginzburg–Landau equations in two dimensions in [23] and the references therein. The authors demonstrate the existence of weakly deformed stripe patterns, which are small perturbations of the uniform vertical stripes, where the small parameter measures the amplitude of the spatial inhomogeneity. Although these studies are for fixed parameters, it might be of interest to study the spatio-temporal dynamics that result from the interactions between these types of heterogeneities and the slow passage of a parameter through a Hopf bifurcation.

Appendix A. Calculation of $A_p(x, y, \mu)$ using the saddle point method

In this appendix, we apply the saddle point method to find the asymptotic formula (2.5) for the particular solution $A_p(x, y, \mu)$ of the linearized CGL PDE. In the complex $\tilde{\mu}$ plane ($\tilde{\mu} = \tilde{\mu}_R + i\tilde{\mu}_I$), the phase function in B_p is

$$\phi + i\psi = -\frac{1}{2}(\tilde{\mu} + i\omega_0)^2,$$

where

$$\phi = -\frac{1}{2}(\tilde{\mu}_R^2 - (\tilde{\mu}_I + \omega_0)^2) \quad \text{and} \quad \psi = -\tilde{\mu}_R(\tilde{\mu}_I + \omega_0).$$

This phase has a saddle point at $\tilde{\mu} = -i\omega_0$. The level sets of ϕ are hyperbolas and also known as Stokes lines. The Stokes lines with $\phi = 0$ through the saddle (which are the asymptotes of the hyperbolas) bound the valleys and hills. They may be parametrized by $\tilde{\mu}_R$ via $\tilde{\mu}_I = \pm\tilde{\mu}_R - \omega_0$. Also, the level sets of ψ are hyperbolas (with asymptotes given by the axes), and they are referred to as anti-Stokes lines.

Let $\delta > 0$ be a small number, independent of ε . We fix an arbitrary value of $\mu \in [\delta, \omega_0]$. We consider the contour $C_r = C_{r1} \cup C_{r2} \cup C_{r3} \cup C_{r4}$, where $C_{r1} = [\mu_0, -\omega_0]$; C_{r2} is the segment of the Stokes line $\tilde{\mu}_I = -\tilde{\mu}_R - \omega_0$ from $-\omega_0$ down to the saddle at $-i\omega_0$; C_{r3} is the segment of the Stokes line $\tilde{\mu}_I = \tilde{\mu}_R - \omega_0$ from the saddle up to the point $q_r = \sqrt{\omega_0\mu} + i(\sqrt{\omega_0\mu} - \omega_0)$, for this fixed value of μ ; and C_{r4} consists of the segment of the steepest ascent curve $\psi = -\omega_0\mu$ from q_r up to the point μ (see Figure A.1).

We evaluate the integrals in the same manner as in Section 2.3 of [17]. We take any initial $B_p(x, y, \mu_0)$ with $\mu_0 \leq -\omega_0$ on or near the attracting QSS, and we track it along C_r to the fixed value μ . At that point,

$$B_p(x, y, \mu) = I_{r1} + I_{r2} + I_{r3} + I_{r4},$$

where, for $\delta \leq \mu \leq \omega_0$,

$$I_{rj} = \frac{1}{\sqrt{\varepsilon}} \int_{C_{rj}} g(x, y, \mu - \tilde{\mu}) e^{-(\tilde{\mu} + i\omega_0)^2/2\varepsilon} d\tilde{\mu}, \quad j = 1, 2, 3, 4.$$

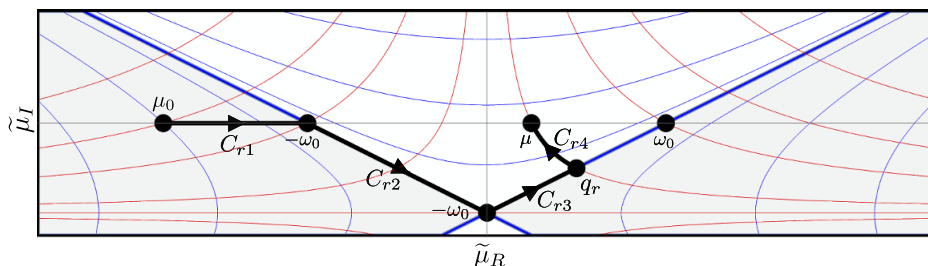


FIGURE A.1. The contour $C_r = C_{r1} \cup C_{r2} \cup C_{r3} \cup C_{r4}$ in the complex $\tilde{\mu}$ plane.

The integral \mathcal{I}_{r1} along the segment $C_{r1} = [\mu_0, -\omega_0]$ is evaluated using the method of steepest descents along the contour formed by the union of the Stokes line through $\mu = -\mu_0$ out to $-\infty - i\omega_0$ and the Stokes line back from $-\infty - i\omega_0$ up to $\mu = -\omega_0$. The dominant contribution comes from the final segment along the latter Stokes line near $\mu = -\omega_0$.

$$\mathcal{I}_{r1} = \left\{ \frac{\sqrt{\varepsilon}}{2\omega_0} (1+i)g(x, y, \mu + \omega_0) + O(\varepsilon^{3/2}) \right\} e^{i\omega_0^2/\varepsilon}. \quad (\text{A.1})$$

Next, we parametrize C_{r2} by $\tilde{\mu}_R$, with $\tilde{\mu}_I(\tilde{\mu}_R) = -(\tilde{\mu}_R + \omega_0)$ and $\tilde{\mu}_R : -\omega_0 \rightarrow 0$. Hence, $-(\tilde{\mu} + i\omega_0)^2/2 = i\tilde{\mu}_R^2$ for all $\tilde{\mu}$ on C_{r2} . It is purely imaginary, corresponding to the fact that C_{r2} lies on a Stokes line $\phi = 0$. Hence, for each $\tilde{\mu}$ on C_{r2} , the integrand is of the form to which the method of stationary phase applies; namely, $g \cdot e^{(i/\varepsilon)h(\tilde{\mu}_R)}$ with $h(\tilde{\mu}_R) = \tilde{\mu}_R^2$. Moreover, the end point $\tilde{\mu} = -i\omega_0$ of C_{r2} ($\tilde{\mu}_R = 0$) is a point of stationary phase, since $h'(0) = 0$ and $h''(0) = 2 \neq 0$, and it is the only such point along C_{r2} . (For the general method in which an end point is a saddle (or turning) point, see, for example, Section 4.1 of [34], in particular formula (4.14).)

Applying the method of stationary phase, we insert the parametrization of C_{r2} , use $\tilde{v} = -\tilde{\mu}_R$, Taylor expand about $\tilde{v} = 0$ (that is, $\tilde{\mu} = -i\omega_0$) and observe that the dominant contribution asymptotically comes from the point of stationary phase at the saddle,

$$\begin{aligned} \mathcal{I}_{r2} &= \frac{1}{\sqrt{\varepsilon}} \int_{-\omega_0}^0 g(x, y, \mu - [\tilde{\mu}_R - i(\tilde{\mu}_R + \omega_0)]) e^{i\tilde{\mu}_R^2/\varepsilon} (1-i) d\tilde{\mu}_R \\ &= \frac{1}{\sqrt{\varepsilon}} (1-i) \int_0^{\omega_0} g(x, y, \mu + [\tilde{v} + i(-\tilde{v} + \omega_0)]) e^{i\tilde{v}^2/\varepsilon} d\tilde{v} \\ &= \frac{1}{\sqrt{\varepsilon}} (1-i) g(x, y, \mu + i\omega_0) \int_0^{\omega_0} (1 + O(\tilde{v})) e^{i\tilde{v}^2/\varepsilon} d\tilde{v} \\ &= \sqrt{\frac{\pi}{2h''(0)}} e^{i\pi/4} (1-i) g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}) \\ &= \sqrt{\frac{\pi}{2}} g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}) \quad \text{for any } \mu \in [\delta, \omega_0]. \end{aligned} \quad (\text{A.2})$$

This leading order term in \mathcal{I}_{r2} will turn out to be half of the leading order term in the total integral for $B_p(x, y, \mu)$ for each $\mu \in [\delta, \omega_0]$.

Next, we show that \mathcal{I}_{r3} gives the other half of the leading order term in B_p . From the definition of C_{r3} , we have that, for any $\mu \in [\delta, \omega_0]$,

$$\mathcal{I}_{r3} = \frac{1}{\sqrt{\varepsilon}} \int_{-i\omega_0}^{q_r} g(x, y, \mu - \tilde{\mu}) e^{-(\tilde{\mu} + i\omega_0)^2/2\varepsilon} d\tilde{\mu}.$$

We use $\tilde{\mu}_R$ to parametrize C_{r3} as $\tilde{\mu}_I = \tilde{\mu}_R - \omega_0$, now with $\tilde{\mu}_R : 0 \rightarrow \sqrt{\omega_0\mu}$. Hence, $-(\tilde{\mu} + i\omega_0)^2/2 = -i\tilde{\mu}_R^2$ along C_{r3} ; and, for each $\tilde{\mu}$ on C_{r3} , the integral is also of the form to apply the method of stationary phase, $g \cdot e^{(i/\varepsilon)\tilde{h}(\tilde{\mu}_R)}$, with $\tilde{h}(\tilde{\mu}_R) = -\tilde{\mu}_R^2$. Moreover, the

initial point $\tilde{\mu} = -i\omega_0$ of C_{r3} ($\tilde{\mu}_R = 0$) is a stationary phase point, since $\tilde{h}'(0) = 0$ and $\tilde{h}''(0) = -2 \neq 0$, and it is the only such point along C_{r3} . We find

$$\begin{aligned} \mathcal{I}_{r3} &= \frac{1}{\sqrt{\varepsilon}} \int_0^{\sqrt{\omega_0\mu}} g(x, y, \mu - [\tilde{\mu}_R + i(\tilde{\mu}_R - \omega_0)]) e^{-i\tilde{\mu}_R^2/\varepsilon} (1+i) d\tilde{\mu}_R \\ &= \frac{1}{\sqrt{\varepsilon}} g(x, y, \mu + i\omega_0) \int_0^{\sqrt{\omega_0\mu}} e^{-i\tilde{\mu}_R^2/\varepsilon} (1 + O(\tilde{\mu}_R)) (1+i) d\tilde{\mu}_R \\ &= \sqrt{\frac{\pi}{2|\tilde{h}''(0)|}} g(x, y, \mu + i\omega_0) e^{-i\pi/4} (1+i) + O(\sqrt{\varepsilon}) \\ &= \sqrt{\frac{\pi}{2}} g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}) \quad \text{for any } \mu \in [\delta, \omega_0]. \end{aligned} \quad (\text{A.3})$$

Finally, we calculate \mathcal{I}_{r4} . Implicitly parametrize C_{r4} using σ ,

$$C_{r4}: \quad -\frac{1}{2}(\tilde{\mu} + i\omega_0)^2 = -\frac{1}{2}(\mu + i\omega_0)^2 + \sigma.$$

The parameter σ starts from $-(\omega_0^2 - \mu^2)/2$ at the point q_r and increases monotonically along C_{r4} to zero at μ . The explicit representation is

$$\tilde{\mu}(\sigma) = -i\omega_0 + [(\mu + i\omega_0)^2 - 2\sigma]^{1/2}.$$

The integration along C_{r4} gives

$$\mathcal{I}_{r4} = \sqrt{\varepsilon} e^{-(\mu+i\omega_0)^2/2\varepsilon} \int_{-(\omega_0^2-\mu^2)/2}^0 \frac{g(x, y, \mu + i\omega_0 - [(\mu + i\omega_0)^2 - 2\sigma]^{1/2}) e^{\sigma/\varepsilon}}{\sqrt{(\mu + i\omega_0)^2 - 2\sigma}} d\sigma.$$

Hence, one finds

$$\begin{aligned} \mathcal{I}_{r4} &= \left[-\frac{\sqrt{\varepsilon} I_a(x)}{\mu + i\omega_0} + \varepsilon^{3/2} \left(\frac{I_a(x, y) + d(\mu + i\omega_0) \Delta I_a(x, y)}{(\mu + i\omega_0)^3} \right) \right. \\ &\quad \left. + O\left(\frac{\varepsilon^{5/2}}{(\mu + i\omega_0)^5} \right) \right] e^{-(\mu+i\omega_0)^2/2\varepsilon}. \end{aligned} \quad (\text{A.4})$$

Summing (A.1), (A.2), (A.3) and (A.4), we have that, for $\delta \leq \mu \leq \omega_0$,

$$\begin{aligned} B_p(x, y, \mu) &= \mathcal{I}_{r1} + \mathcal{I}_{r2} + \mathcal{I}_{r3} + \mathcal{I}_{r4} \\ &= -\frac{\sqrt{\varepsilon} I_a(x, y)}{\mu + i\omega_0} e^{-(\mu+i\omega_0)^2/2\varepsilon} \\ &\quad + \varepsilon^{3/2} \left(\frac{I_a(x, y) + d(\mu + i\omega_0) \Delta I_a(x, y)}{(\mu + i\omega_0)^3} \right) e^{-(\mu+i\omega_0)^2/2\varepsilon} \\ &\quad + O\left(\frac{\varepsilon^{5/2} e^{-(\mu+i\omega_0)^2/2\varepsilon}}{(\mu + i\omega_0)^5} \right) + \sqrt{2\pi} g(x, y, \mu + i\omega_0) + O(\sqrt{\varepsilon}). \end{aligned}$$

Finally, we translate the formula back to the dependent variable A , which completes the derivation of (2.5).

Acknowledgements

The authors congratulate Bernd Krauskopf on the occasion of this sixtieth birthday. We are pleased to help honour his many contributions to the fields of bifurcation theory, dynamical systems, delay differential equations, numerical methods for dynamical systems and applications of dynamical systems.

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