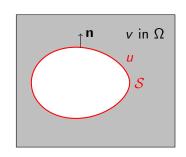
### Closest Point Method: coupled surface-bulk problems

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Oriel CBL course, Oxford, Aug 2014

# Surface bulk coupling: diffusion problem (1)



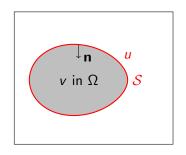
$$u_t = D\Delta_{\mathcal{S}} u - \mu u + \nu v|_{\mathcal{S}}, \text{ on } \mathcal{S}$$
 (1a)  
$$v_t = d\Delta v, \text{ in } \Omega$$
 (1b)

Boundary condition(B.C.) for v on S:

$$d\frac{\partial \mathbf{v}}{\partial \mathbf{n}}\bigg|_{\mathcal{S}} = \mu \mathbf{u} - \nu \mathbf{v}|_{\mathcal{S}}.$$
 (1c)

Prescribed B.C.s on four edges of the box (depending on physics and what we care).

# Surface bulk coupling: diffusion problem (2)

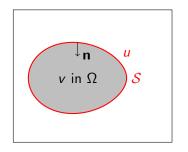


$$\mathbf{u}_t = D\Delta_{\mathcal{S}}\mathbf{u} - \mu\mathbf{u} + \nu\mathbf{v}|_{\mathcal{S}}, \text{ on } \mathcal{S}$$
 (2a)  
 $\mathbf{v}_t = d\Delta\mathbf{v}, \text{ in } \Omega$  (2b)

Boundary condition(B.C.) for v on S:

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# Surface bulk coupling: diffusion problem (2)



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 (2c)

For convenience of illustration, we will focus on problem (2).

## Big picture: forward Euler alternating time-stepping

Starting from  $\mathbf{u}^0$  and  $\mathbf{v}^0$  satisfying  $d\frac{\partial \mathbf{v}^0}{\partial \mathbf{n}}\big|_{\mathcal{S}} = \mu \mathbf{u}^0 - \nu \mathbf{v}^0|_{\mathcal{S}}$ , for  $n=0,1,2,\cdots$ , alternating between the following two steps:

(1) Taking the bulk function v as known, evolve the surface function u:

$$u^{n+1} = u^n + \tau (D\Delta_{\mathcal{S}} u^n - \mu u^n + \nu v^n|_{\mathcal{S}}).$$

(2) Taking the surface function u as known, evolve the bulk function v:

$$v^{n+1} = v^n + \tau d\Delta v^n$$
 with B.C.  $d\frac{\partial v^{n+1}}{\partial \mathbf{n}}\bigg|_{\mathcal{S}} = \mu u^{n+1} - \nu v^{n+1}|_{\mathcal{S}}.$ 

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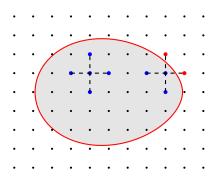
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 with B.C.  $d\frac{\partial v^{n+1}}{\partial \mathbf{n}}\bigg|_{\mathcal{S}} = \mu u^{n+1} - \nu v^{n+1}|_{\mathcal{S}}.$ 

We already know how to do (1) with the Closest Point Method, now we focus on step (2).



- The Laplacian stencil of some grid points might have points outside the computational domain  $\Omega$ . How do we deal with these grid points?
- How to impose the boundary condition  $d \frac{\partial v}{\partial \mathbf{n}} |_{\mathcal{S}} = \mu \mathbf{u} \nu v |_{\mathcal{S}}$ ?

### Approximation of boundary condition

According to the boundary condition  $d\frac{\partial v}{\partial \mathbf{n}}|_{\mathcal{S}} = \mu \mathbf{u} - \nu \mathbf{v}|_{\mathcal{S}}$ , for grid point  $\mathbf{x}_i$  outside domain  $\Omega$ , we have the following approximation:

$$-d\frac{v(\mathbf{x}_i) - v(\bar{\mathbf{cp}}(\mathbf{x}_i))}{\|\mathbf{x}_i - \bar{\mathbf{cp}}(\mathbf{x}_i)\|} = \mu u(\mathbf{cp}(\mathbf{x}_i)) - \nu v(\mathbf{cp}(\mathbf{x}_i)). \tag{3}$$

Here  $\bar{cp}(\mathbf{x}_i) = 2cp(\mathbf{x}_i) - \mathbf{x}_i$  is the mirror image of  $\mathbf{x}_i$  with respect to  $cp(\mathbf{x}_i)$ .

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Here  $c\bar{p}(\mathbf{x}_i) = 2cp(\mathbf{x}_i) - \mathbf{x}_i$  is the mirror image of  $\mathbf{x}_i$  with respect to  $cp(\mathbf{x}_i)$ .

Since  $\operatorname{cp}(\mathbf{x}_i)$  and  $\operatorname{cp}(\mathbf{x}_i)$  are not grid points, we approximate their values by Lagrange interpolation. Denote the interpolation matrix for  $v(\operatorname{cp}(\mathbf{x}_i))$  by  $\mathbf{E}$ , denote the interpolation matrix for  $v(\operatorname{cp}(\mathbf{x}_i))$  by  $\mathbf{E}$ , and denote the distance vector consisting  $\|\mathbf{x}_i - \operatorname{cp}(\mathbf{x}_i)\|$  by  $\operatorname{dist}$ . Assuming  $\mathbf{u}$  is known on the surface, equation (3) can be approximated by:

$$-d(\mathbf{v}-\bar{\mathbf{E}}\mathbf{v})./\mathbf{dist}=\mu\mathbf{u}-\nu\mathbf{E}\mathbf{v}.\tag{4}$$

#### How to evolve **v**?

Our goal is:

$$v^{n+1} = v^n + \tau d\Delta v^n$$
 with B.C.  $d\frac{\partial v^{n+1}}{\partial \mathbf{n}}\bigg|_{\mathcal{S}} = \mu u^{n+1} - \nu v^{n+1}|_{\mathcal{S}}.$ 

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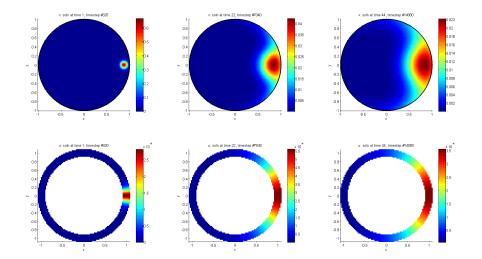
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Step 1: evolve by the normal Laplacian matrix for all grid points.

$$\mathbf{v}^{n+1} = \mathbf{v}^n + \tau d \mathbf{L} \mathbf{v}^n$$

Step 2: correct for grid points outside Ω to match the B.C..
Let Band be the set of computational grid points covering Ω, innerBand be the subset of Band containing grid points inside Ω or on S, and outerBand be the subset of Band containing grid points outside Ω. The solution vector v can be written as v = [v(innerBand), v(outerBand)]<sup>T</sup>. From equation (4), taking v(innerBand) as known, solve for v(outerBand).

## Surface-bulk coupling: numerical results

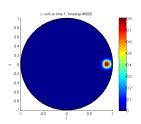


# Surface bulk coupling: Fisher-KPP equation in the bulk

#### (Dawson 2014)

$$egin{array}{lll} u_t &=& D\Delta_{\mathcal{S}} u - \mu u + 
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u v, \end{array}$$

 $\mathbf{n}$ : unit normal pointing inwards  $\Omega$ .

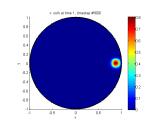


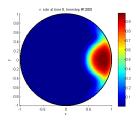
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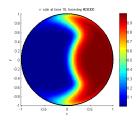
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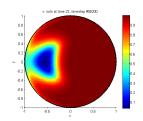
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 $\mathbf{n}$ : unit normal pointing inwards  $\Omega$ .









## Surface bulkcoupling: reaction-diffusion process

(Macdonald, Merriman, Ruuth, 2013)

$$u_t = f(u, v) + D_u \Delta_S u - \alpha_1 u + \beta_1 U$$
, on  $S$ 

$$v_t = g(u, v) + D_v \Delta_S v - \alpha_2 v + \beta_2 V$$
, on  $S$ 

$$U_t = f(U, V) + D_U \Delta U$$
, in  $\Omega$ 

$$V_t = g(U, V) + D_V \Delta V$$
, in  $\Omega$ 

with coupling boundary conditions:

$$D_U \frac{\partial U}{\partial n} = \alpha_1 u - \beta_1 U, \ D_V \frac{\partial V}{\partial n} = \alpha_2 v - \beta_2 V$$

