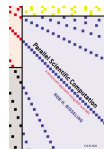
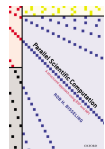


# ***Laplacian Matrices*** **(PSC §4.8)**



# Physical domain

- In many applications, a physical domain exists that can be distributed naturally by assigning a **contiguous subdomain** to every processor.
- Communication is only needed for exchanging information across the **subdomain boundaries**.
- Often, the domain is structured as a **multidimensional rectangular grid**, where grid points interact only with a set of immediate neighbours.
- In the 2D case, these could be the neighbours to the north, east, south, and west.
- An example is the **heat equation**, where the value at a grid point represents the temperature at the corresponding location.



# 2D Laplacian operator for $k \times k$ grid

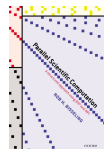
Compute

$$\Delta_{i,j} = x_{i-1,j} + x_{i+1,j} + x_{i,j+1} + x_{i,j-1} - 4x_{i,j}, \quad \text{for } 0 \leq i, j < k,$$

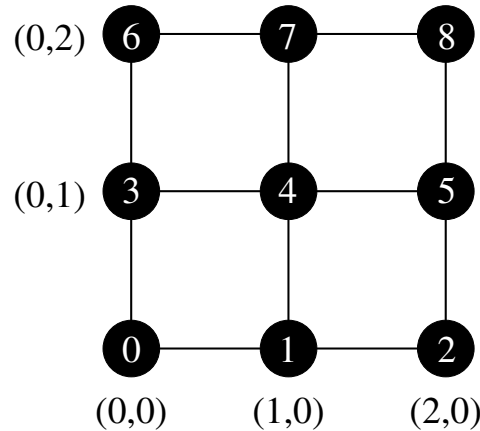
where  $x_{i,j}$  denotes the temperature at grid point  $(i, j)$ .

By convention,  $x_{i,j} = 0$  outside the grid.

- $x_{i+1,j} - x_{i,j}$  approximates the **derivative** of the temperature in the  $i$ -direction.
- $(x_{i+1,j} - x_{i,j}) - (x_{i,j} - x_{i-1,j}) = x_{i-1,j} + x_{i+1,j} - 2x_{i,j}$  approximates the **second derivative**.



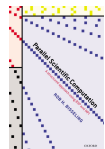
# Relation grid–vector



- A  $3 \times 3$  **grid**, which corresponds to a **vector** of length 9. For each grid point  $(i, j)$ , the index  $i + 3j$  of the corresponding vector component is shown.
- More in general,

$$v_{i+jk} \equiv x_{i,j}, \quad u_{i+jk} \equiv \Delta_{i,j},$$

for  $0 \leq i, j < k$ .

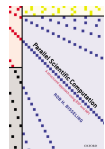


# Relation operator–matrix

$$A = \begin{bmatrix} -4 & 1 & \cdot & 1 & \cdot & \cdot & \cdot & \cdot & \cdot \\ 1 & -4 & 1 & \cdot & 1 & \cdot & \cdot & \cdot & \cdot \\ \cdot & 1 & -4 & \cdot & \cdot & 1 & \cdot & \cdot & \cdot \\ 1 & \cdot & \cdot & -4 & 1 & \cdot & 1 & \cdot & \cdot \\ \cdot & 1 & \cdot & 1 & -4 & 1 & \cdot & 1 & \cdot \\ \cdot & \cdot & 1 & \cdot & 1 & -4 & \cdot & \cdot & 1 \\ \cdot & \cdot & \cdot & 1 & \cdot & \cdot & -4 & 1 & \cdot \\ \cdot & \cdot & \cdot & \cdot & 1 & \cdot & 1 & -4 & 1 \\ \cdot & \cdot & \cdot & \cdot & \cdot & 1 & \cdot & 1 & -4 \end{bmatrix}$$

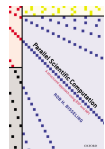
$$\mathbf{u} = A\mathbf{v} \iff$$

$$\Delta_{i,j} = x_{i-1,j} + x_{i+1,j} + x_{i,j+1} + x_{i,j-1} - 4x_{i,j}, \quad \text{for } 0 \leq i, j < k.$$



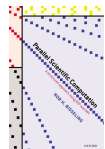
## *Domain view vs. matrix view*

- In general, it is best to view the Laplacian as an operator on the physical domain.
- This **domain view** has the advantage that it naturally leads to the use of a regular data structure.
- Occasionally, however, it may be beneficial to **view** the Laplacian as a **matrix**, so that we can apply our knowledge about sparse matrix–vector multiplication.

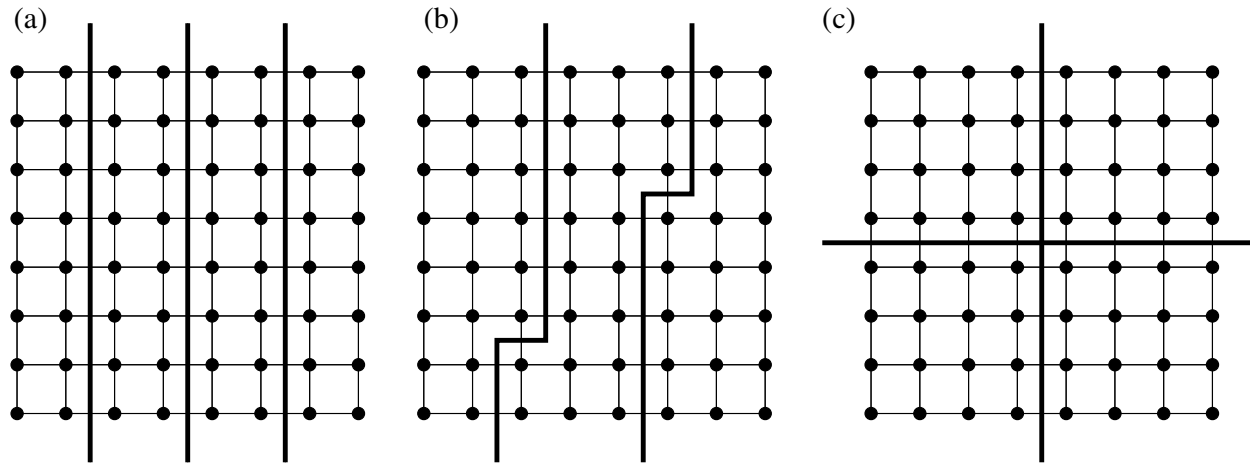


## Find a domain distribution

- Here, we adopt the domain view, so that we must assign each grid point to a processor.
- We assign the values  $x_{i,j}$  and  $\Delta_{i,j}$  to the owner of grid point  $(i, j)$ , which translates into  $\text{distr}(\mathbf{u}) = \text{distr}(\mathbf{v})$ .
- We use a row distribution for the matrix and assign row  $i + jk$  to the same processor as vector component  $u_{i+jk}$  and hence grid point  $(i, j)$ .
- The resulting sparse matrix–vector multiplication algorithm has **two supersteps**, the fanout and the local matrix–vector multiplication.
- The computation time for an **interior** point is 5 flops; for a **border** point 4 flops; for a **corner** point 3 flops.



# Distribution into strips and blocks

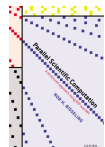


- (a) Distribution into strips: long Norwegian borders,

$$T_{\text{comm, strips}} = 2kg.$$

- (b) Boundary corrections improve load balance.
- (c) Distribution into square blocks: shorter borders,

$$T_{\text{comm, squares}} = \frac{4k}{\sqrt{p}}g \quad (\text{for } p > 4).$$



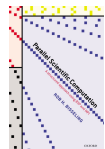


# Surface-to-volume ratio

- The communication-to-computation ratio for square blocks is

$$\frac{T_{\text{comm, squares}}}{T_{\text{comp, squares}}} = \frac{4k/\sqrt{p}}{5k^2/p} g = \frac{4\sqrt{p}}{5k} g.$$

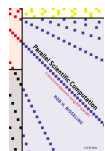
- This ratio is often called the surface-to-volume ratio, because in 3D the surface of a domain represents the communication with other processors and the volume represents the amount of computation of a processor.



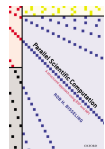
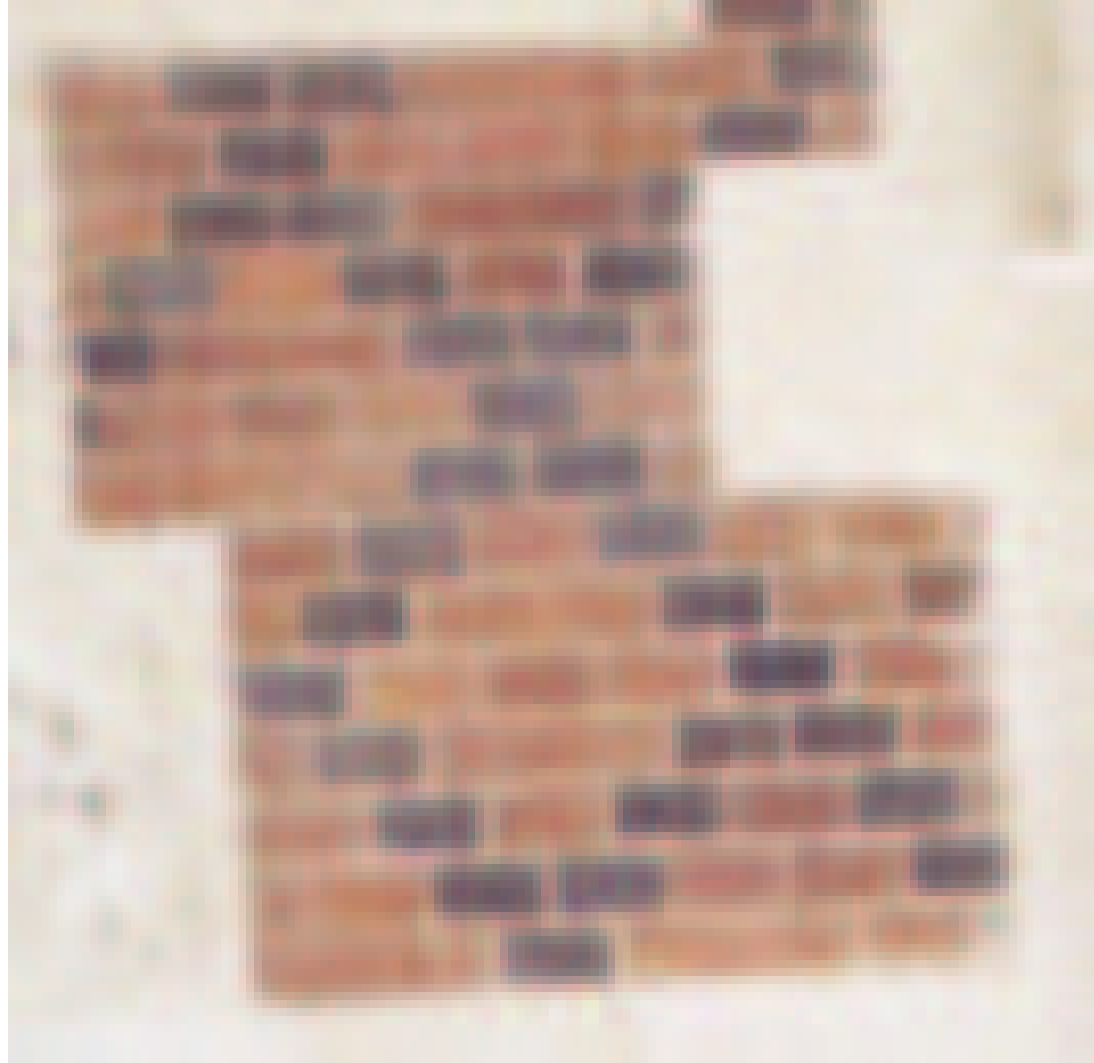
# *What do we do at scientific workshops?*



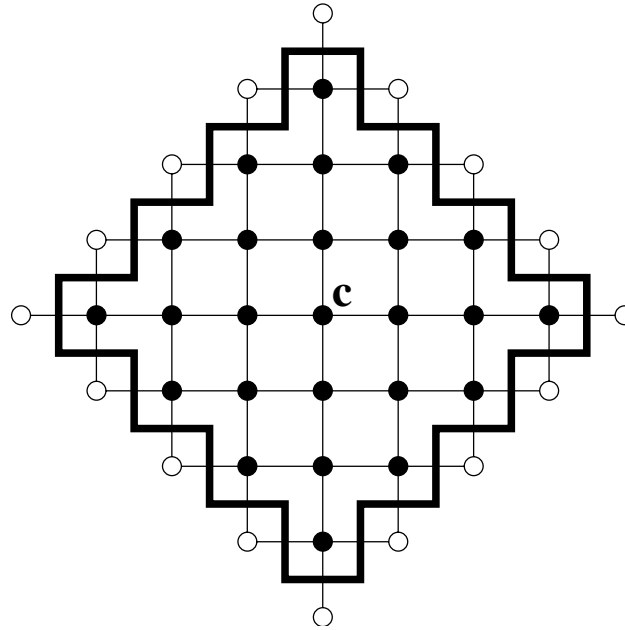
Participants of HLPP 2001, International Workshop on **High-Level** Parallel Programming, Orléans, France, June 2001, studying Château de Blois.



# *The high-level object of our study*



## Blocks are nice, diamonds ...



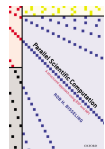
$$r = 3$$

- Digital diamond, or closed  $l_1$ -sphere, defined by

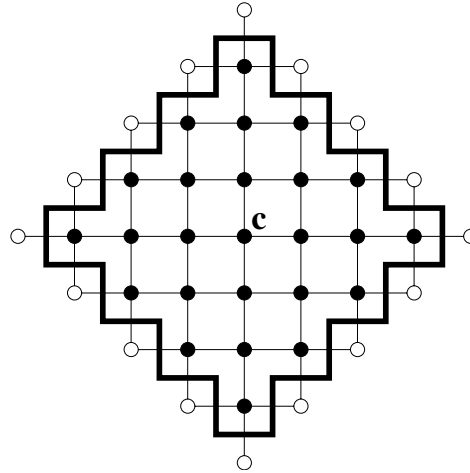
$$B_r(c_0, c_1) = \{(i, j) \in \mathbf{Z}^2 : |i - c_0| + |j - c_1| \leq r\},$$

for integer radius  $r \geq 0$  and centre  $\mathbf{c} = (c_0, c_1) \in \mathbf{Z}^2$ .

- $B_r(\mathbf{c})$  is the set of points with Manhattan distance  $\leq r$  to the central point  $\mathbf{c}$ .



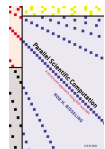
# Points of a diamond



$$r = 3$$

The number of points of  $B_r(c)$  is

$$\begin{aligned}
 & 1 + 3 + 5 + \cdots + (2r - 1) + (2r + 1) + (2r - 1) + \cdots + 1 \\
 &= 2 \sum_{k=0}^{r-1} (2k + 1) + (2r + 1) = 4 \sum_{k=0}^{r-1} k + 4r + 1 \\
 &= 2(r - 1)r + 4r + 1 = 2r^2 + 2r + 1.
 \end{aligned}$$



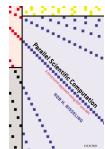
The number of neighbouring points is  $4r + 4$ .

# *Diamonds are forever*

- Assume that the diamond has its fair share  $2r^2 + 2r + 1 = \frac{k^2}{p}$  of the grid points.
- Therefore,  $2r^2 \approx \frac{k^2}{p}$  for large  $r$ , and hence  $r \approx \frac{k}{\sqrt{2p}}$ .
- Just on the basis of  $4r + 4$  receives, we have

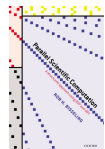
$$\frac{T_{\text{comm, diamonds}}}{T_{\text{comp, diamonds}}} = \frac{4r + 4}{5(2r^2 + 2r + 1)}g \approx \frac{2}{5r}g \approx \frac{2\sqrt{2p}}{5k}g.$$

- Compare with value  $\frac{4\sqrt{p}}{5k}g$  for square blocks:  
factor  $\sqrt{2}$  less.
- Gain caused by reuse of data: value at grid point is used twice but sent only once.

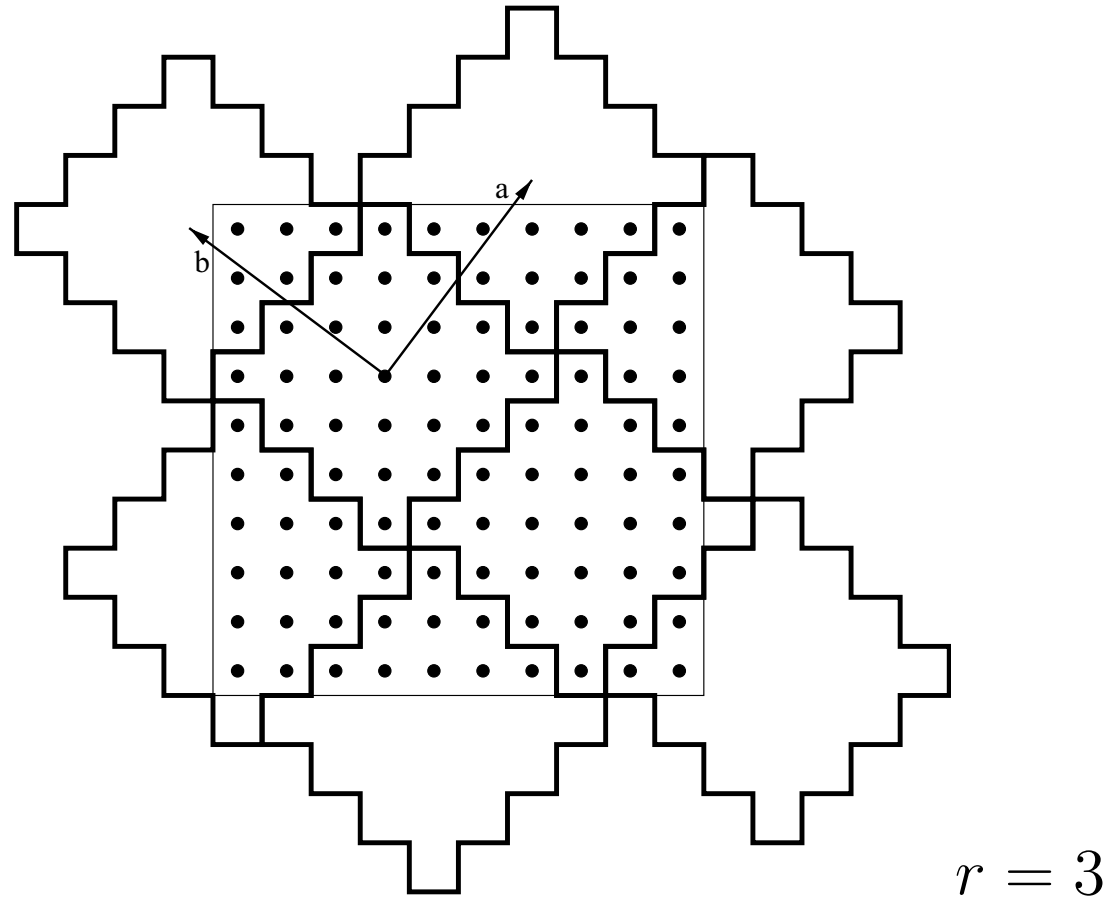




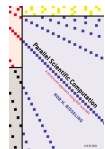
# *Alhambra: tile the whole space*



# *Tile the whole sky with diamonds*



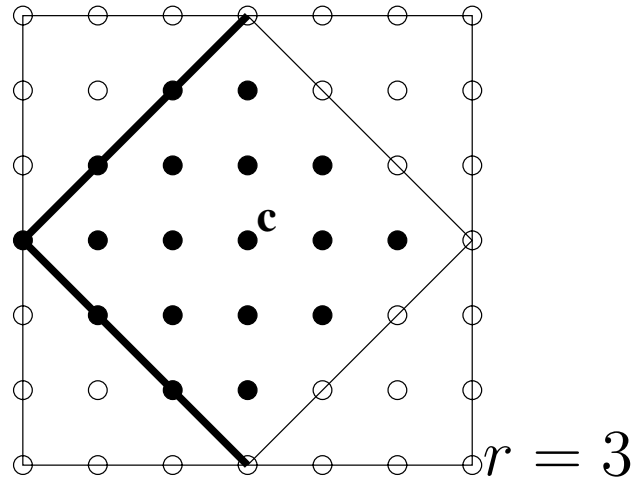
Diamond centres at  $\mathbf{c} = \lambda \mathbf{a} + \mu \mathbf{b}$ ,  $\lambda, \mu \in \mathbf{Z}$ ,  
where  $\mathbf{a} = (r, r + 1)$  and  $\mathbf{b} = (-r - 1, r)$ .



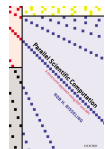
Good method for an **infinite grid**.



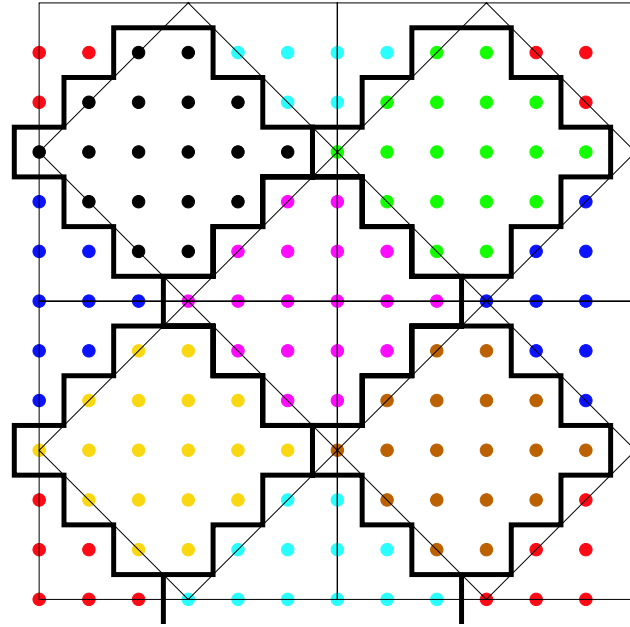
# Practical method for finite grids



- Discard one layer of points from the north-eastern and south-eastern border of the diamond.
- For  $r = 3$ , the number of points decreases from 25 to 18.

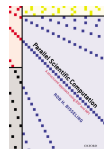


# $12 \times 12$ *computational grid: periodic partitioning*

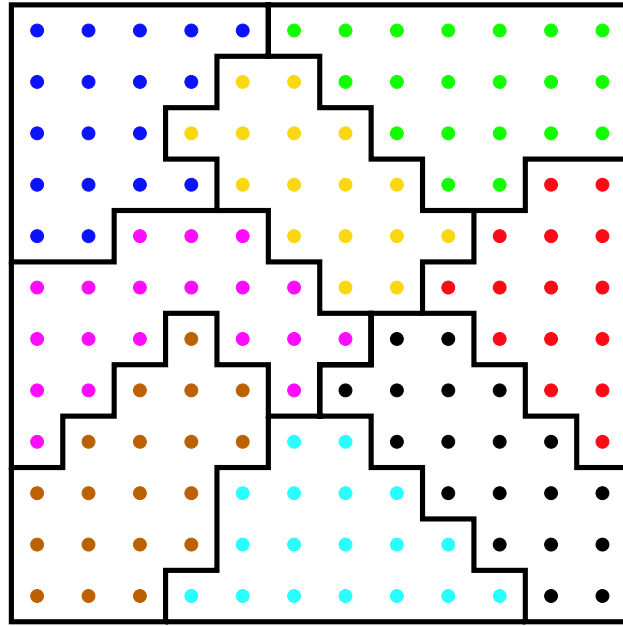


8 processors

- Total computation: 672 flops. Avg 84. Max 90.
- Communication: 104 values. Avg 13. Max 14.
- Total time:  $90 + 14g = 90 + 14 \cdot 10 = 230$  (ignoring  $2l$ ).
- Rectangular  $6 \times 3$  blocks: time would be  $87 + 15 \cdot 10 = 237$ . Worse!

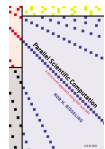


# $12 \times 12$ computational grid: Mondriaan partitioning



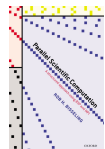
8 processors

- Total computation: 672 flops. Avg 84. Max 91. ( $\epsilon = 10\%$ .)
- Communication: 85 values. Avg 10.525. Max 16.
- Total time:  $91 + 16g = 91 + 16 \cdot 10 = 251$ .
- **Challenge:** better solution can be obtained manually, using ideas from both solutions shown. Current best known solution is 200 (Vermolen 2005).

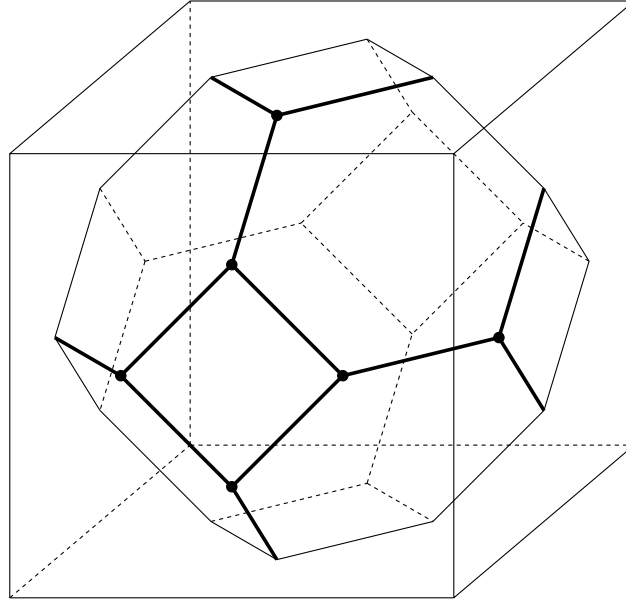


## Three dimensions

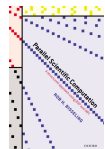
- If a processor has a cubic block of  $N = k^3/p$  points, about  $\frac{6k^2}{p^{2/3}} = 6N^{2/3}$  are boundary points. In 2D, only  $4N^{1/2}$ .
- If a processor has a  $10 \times 10 \times 10$  block, 488 points are on the boundary. About half!
- Thus, communication is important in 3D.
- Based on the surface-to-volume ratio of a 3D digital diamond, we can aim for a reduction by a factor  $\sqrt{3} \approx 1.73$  in communication cost.
- The prime application of diamond-shaped distributions will most likely be in 3D.



# Basic cell for 3D



- Basic cell: grid points in a **truncated octahedron**.
- For load balancing, take care with the boundaries.
- What You See, Is What You Get (WYSIWYG):  
4 hexagons and 3 squares visible at the front are included. Also 12 edges, 6 vertices.
- Gain factor of 1.68 achieved for  $p = 2q^3$ .

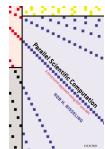


# Comparing 3 distribution methods in 2D and 3D

Grid	$p$	Rectangular	Diamond	Mondriaan
$1024 \times 1024$	2	1024	2046	1024
	4	1024	2048	1240
	8	1280	1026	1378
	16	1024	1024	1044
	32	768	514	766
	64	512	512	548
	128	384	258	395
$64 \times 64 \times 64$	16	4096	2402	2836
	128	1024	626	829

Communication cost (in  $g$ ) for a Laplacian operation on a grid.

Mondriaan with  $\epsilon = 10\%$ .



# Summary

- Communication can be reduced tremendously by using **knowledge of the physical domain**.
- To achieve a good distribution with a low surface-to-volume ratio, **all dimensions must be cut**. In 2D, this gives square blocks. In 3D, cubic subdomains.
- In 2D, an even better method is to use **digital diamonds** (with two edge layers removed). This basic cell can be used to tile a rectangular domain in a straightforward manner. Best performance is obtained for  $p = 2q^2$ .
- In 3D, the best method is to use **truncated octahedra** with WYSIWYG tie breaking at the boundaries. Best performance is obtained for  $p = 2q^3$ .
- In 3D, the performance of Mondriaan is between that of cubes and truncated octahedra.

