TD11: Discrete Capacity-Constrained Voronoi Tessellation

In this TP, the objective is to implement a point sampling strategy from a variant of Voronoi diagram: the capacity-constrained Voronoi diagram \[1, 2\].

The main idea is to optimize a Voronoi diagram (by moving sites) such that each Voronoi cell have the same area (capacity). The digital version can be described as follows: we want to sample \( n \) points in a \( N \times N \) domain. Hence, each cell will be represented by a label \( l \in \{1 \ldots n\} \). The problem is thus to distribute labels to pixels in the \( N \times N \) domain.

The first version of the algorithm can be sketched as follows:

1. First, each pixel randomly picks its label \( l \) such that the number of pixels with label \( l \) is \( \frac{N^2}{n} \);
2. To each label \( l \), \( C_l \) denotes the centroid of points with label \( l \);
3. While stability
   
   - Randomly select a pair of points \( p \) and \( q \) with labels \( l_p \) and \( l_q \) \( (l_p \neq l_q) \);
   - If swapping the labels of \( p \) and \( q \) decreases the energy, we perform the swap \( (l_p \leftrightarrow l_q) \) and update the centroids \( C_p \) and \( C_q \);

The global energy to minimize is related to a “Voronoi-like” energy. In this simple digital setting, we just need to check is swapping the label reduces the distances between the points and their new cluster centroids. More formally, we define

\[
\Delta e(p, l_p, l_q) = ||p - l_p||^2 - ||p - l_q||^2
\]

as the energy difference when the change the label of \( p \) from \( l_p \) to \( l_q \). We perform the swap if both \( \Delta e(p, l_p, l_q) \) and \( \Delta e(q, l_q, l_p) \) are negative.

When no more swap reduces the energy (or when the random pair selection fails to find a good pair for long time), we stop. Doing so, we converge to a structure where labels induce a Voronoi map with Voronoi cells with same area. Here you have an illustration with three samples:

![Illustration](image.png)

**Exercise 1 Naive Algorithm**

Use DGtal and Image data structure of DGtal to implement the naive algorithm as described above (see TP8).

Please consider a centroid data structure to store the position of the \( n \) centroids (maybe a `std::vector<RealPoint>`).

Start to experiment this algorithm on limited domains first.
Exercise 2 Go Optimizations!

In the description of the naive version, the random selection of a pair of points/labels to swap. A first optimization can be described as follows: For each label, we sort the points with respect to their distance to its centroid (decreasing order).

Hence, we use this ordering to first try to swap points which are farthest from their centroid. The argument is that farthest points have an higher probability to be in the wrong “cluster”.

This optimization would allow you to have faster convergence and thus to be able to consider larger domains. In the authors’ original implementation, several further tricks can be used, can you imagine additional optimization heuristics.

References
