

Improving Data-Parallel Construction of Dn-Nets with Maximum Dispersion

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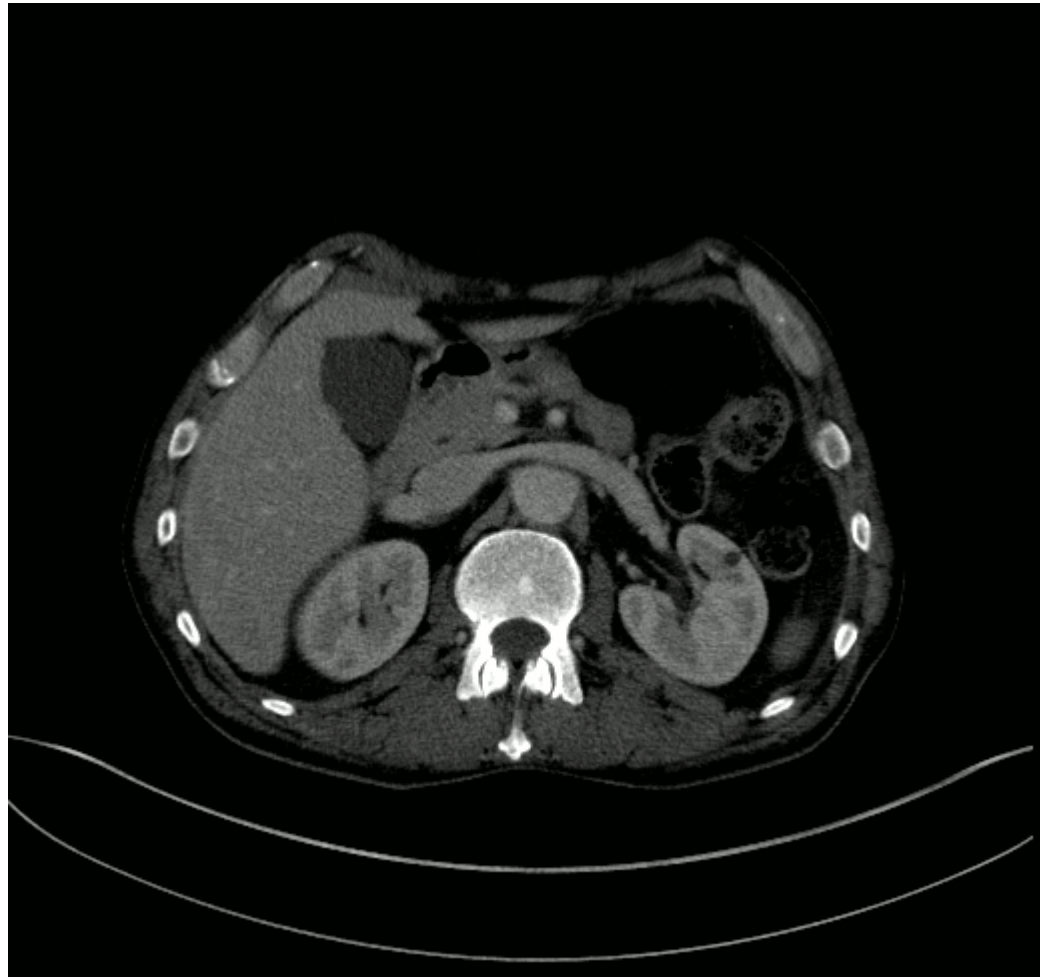
Data-parallel construction of Dn-Nets

Motivation

Joint project: SALK (Salzburger Uniklinik)
Computed Tomography (CAT scans)

Task: Image Segmentation
-> detection of cysts in the body

Sample CAT-scan:



~ 500 patients , 200 - 300 slices
per patient
Trainings set: ~ 140.000 slices

Ground truth:
Kidneys + cysts marked
manually

CAT scan segmentation

Principal Component Analysis

algorithmic creation of 64x64 pixel-patches

auto-correlation matrix

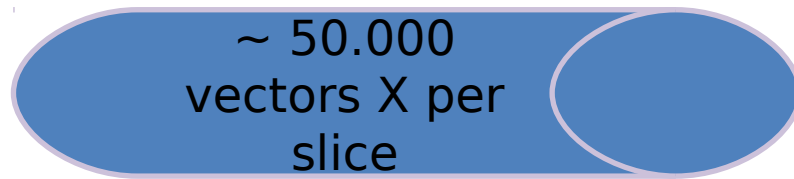
Eigensystem -> preserve 32 most eminent vectors

Code book

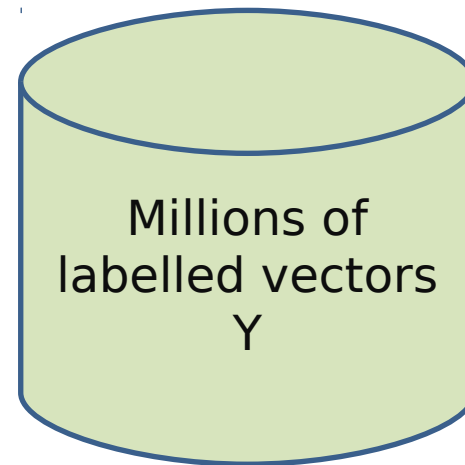
10 – 20 mio. Entries (vectors of 32 floats+labelling data)

Image Segmentation:

Convolution of new patient's CAT scan
book vectors



Code



Task: For each new patch X find nearest neighbor in Y and assign
corresponding label Y_label.

Speeding up the Process:

Options:

Parallel and distributed computing (clusters)

Novel hardware -> GPGPU, CELL BE,
FPGA

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Reduction of code book size -> reduce workload per (kD-trees, sorting, etc. does not work -> curse of dimensionality)

Dn-Nets with Maximum Dispersion

Mathematical concept of Dn-Nets

Metric Dispersion

$$d(x, y) = \left(\sum_{k=1}^s |x_k - y_k|^2 \right)^{1/2}$$

$$\delta_N = \delta(x_1, x_2, \dots, x_N) = \max_{x \in E} (\min(d(x, x_k)))$$

Dn-Net over finite metric space E
approximates high-dimensional data up
to some arbitrary error e

Simpler interpretation: Dn is the radius of
the largest hyper-sphere, that fits within

(x1,x2,... xn).

Construction of Dn-Nets

Choose two seed points (x_1, x_2) with $d(x_1, x_2) = \max (d(dx, dy))$ for all x, y in E

Iteratively add point x , such that $\max(\min(d(x, x_k), k=1, \dots, N))$ is reached.

In other words: at each step we add that point x to our Dn-Net, which decreases the value of D_n by the smallest possible amount.

Previous work

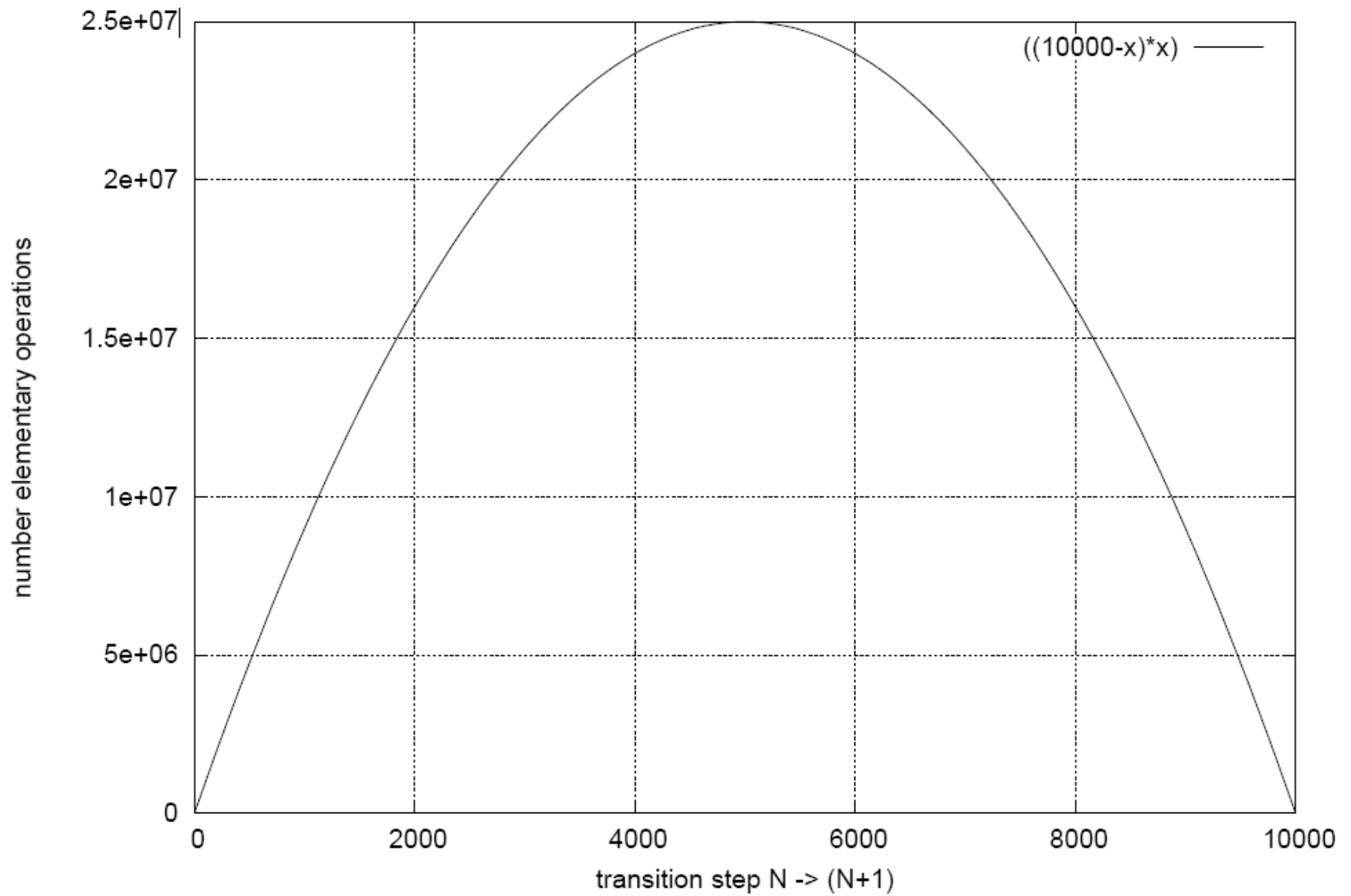
High computational complexity

Transition from $D(n) \rightarrow D(n+1)$: $(E-N)N$
elementary operations

Total complexity: $O(E N^2 - N^3)$

Elementary operation: euclidean distance

distribution of computational complexity

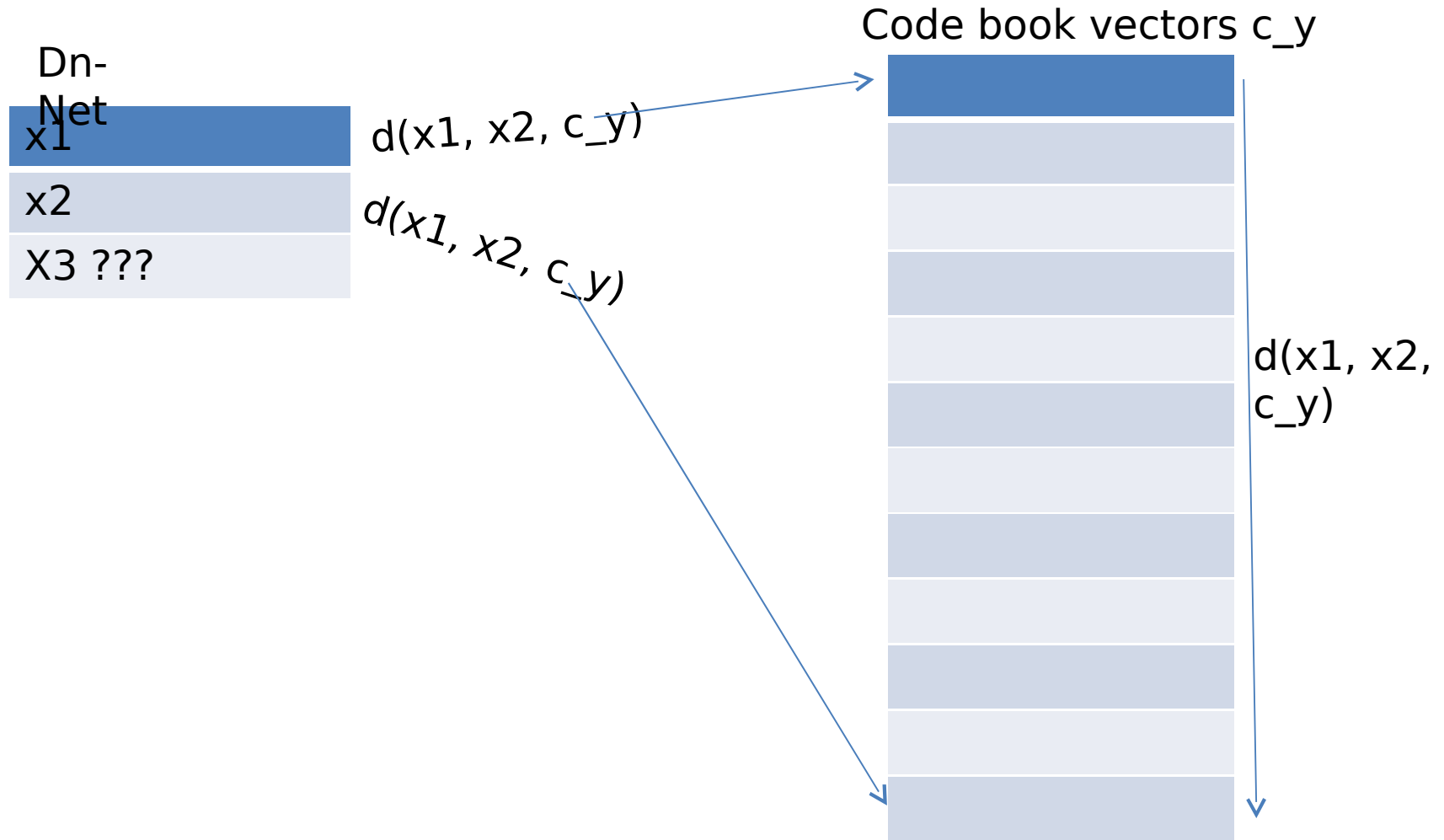


Improved construction of Dn-Nets

Basic idea: instead of (re)computing distances between potential follow-up vectors x_k for a given Dn-Net, the minimum distances for each vector in relation to the Dn-Net are kept in memory.

Also: in the previous algorithm the order in which vectors were added to the Dn-Net had been omitted -> we had to compute distances ($d(DN, x_k)$) over and over again for each new vector.

Previous vs. Improved Algorithm I



Previous vs. Improved Algorithm II

Dn-
Net
x1
x2
x3 ???

We KNOW x2 already !!!
(because it's the second of
our two seed points)

Distances array D holds the current minimum distances
between Dn-Net and entries c1...cN. The only vector that
can change the whole picture is the follow-up-vector x2.
Thus we only have to compare x2 and the values of
d(x1, cx), no matter how big Dn has already grown.

Distances array D

d(x1, c1)
d(x1, c2)
d(x1, c3)
d(x1, c4)
....
d(x1, c12345678)

Previous vs. Improved Algorithm III

x1
x2
x3 ???

Adding x2 involves a simple operation:

For all entries x of the codebook:

Step 1: $C_x = \min(d(x1, c1), d(x2, c1))$ with $d(x1, c1)$ already stored in the array.

Step 2: find $\max(C_x)$,on-the-fly' during the updates, which denotes the next follow-up vector $x_{(n+1)}$. We can iteratively feed this vector into step 1 again and find all follow-up vectors in linear overall complexity.

d(x1, c1)
d(x1, c2)
d(x1, c3)
d(x1, c4)
....
d(x1, c12345678)



Data-Parallelism

al data-independance during update of Distances array D

Computations very well suited for GPGPUs: e.g. CUDA, OpenCL

Implementation by means of two kernels:

- Find seed points x_1, x_2
 - brute force: for each C_n find $\max(d(C_n, C_x))$, $x=(1..\text{codebooksize})$
 - real world data: $\sim 10 - 30$ petaflops needed
- Update Distances D

Results

Computation of some small Dn-Nets on i7 and GPGPU

	i7-CPU, 2.66 GHz	NVIDIA GTX680	NVIDIA GTX 480
Δ_N -net:50k vectors, E=100k vectors	543.94 1.0x	32.70 16.6x	51.75 10.5x
Δ_N -net :500k vectors, E=1m vectors	54433.2 1.0x	2642.84 20.5x	2856.46 19.0x
Δ_N -net:5500 vectors, E=6.3 mio. vectors	7852.6 1.0x	185.67 42.2x	243.97 32.1x
Δ_N -net:15000 vectors, E=6.3 mio. vectors	21392.8 1.0x	499.56 42.8x	652.86 32.7x

Conclusions

presented an improved algorithm for the construction of Dn-Nets, which results in a total computational complexity of $O(E n^2)$ and linear complexity $O(E)$ in each step.

The new algorithm can be considered a big improvement as it enables the construction of Dn-Nets that are common in real-world problems, and which could not be tackled by previous methods even on large clusters of GPGPU-systems.

The presented algorithm lends itself well for GPGPU-execution, which typically enables substantial speedups in the range of 30-40x against single cpu cores and 10x against multi-core CPUs.

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Thank You !

