Các phép biến đổi unitary

Unitary Transforms

Unitary Transforms

- Sort samples *f(x,y)* in an *MxN* image (or a rectangular block in the image) into colunm vector of length *MN*
- Compute transform coefficients

$$\vec{c} = A\vec{f}$$

  where A is a matrix of size *MNxMN*
- The transform A is unitary, iff $A^{-1} = \underbrace{A^{*T} \equiv A^{H}}_{\text{Hermitian conjugate}}$
- If A is real-valued, i.e., $A^{-1}=A^{*}$, transform is „orthonormal“

Energy conservation with unitary transforms

- For any unitary transform $\vec{c} = A\vec{f}$ we obtain

$$\left\|\vec{c}\right\|^2 = \vec{c}^{\,H}\vec{c} = \vec{f}^{\,H}A^{H}A\vec{f} = \left\|\vec{f}\right\|^2$$

- Interpretation: every unitary transform is simply a rotation of the coordinate system.
- Vector lengths („energies“) are conserved.

Energy distribution for unitary transforms

- Energy is conserved, but often will be unevenly distributed among coefficients.
- Autocorrelation matrix

$$R_{cc} = E\left[\vec{c}\vec{c}^{\,H}\right] = E\left[A\vec{f}\cdot\vec{f}^{\,H}A^{H}\right] = AR_{ff}A^{H}$$

- Mean squared values („average energies“) of the coefficients c_i are on the diagonal of R_{cc}

$$E\left[c_i^2\right] = \left[R_{cc}\right]_{i,i} = \left[AR_{ff}A^{H}\right]_{i,i}$$

Eigenmatrix of the autocorrelation matrix

Definition: eigenmatrix Φ of autocorrelation matrix R_{ff}

- Φ is unitary
- The columns of Φ form an orthonormalized set of eigenvectors of R_{ff}, i.e.,

$$R_{ff}\Phi = \Phi\Lambda$$

$$\Lambda = \begin{pmatrix} \lambda_0 & & & 0 \\ & \lambda_1 & & \\ & & \ddots & \\ 0 & & & \lambda_{MN-1} \end{pmatrix}$$

is a diagonal matrix of eigenvalues.

- R_{ff} is symmetric nonnegative definite, hence $\lambda_i \geq 0$ for all i
- R_{ff} is normal matrix, i.e., $R_{ff}^H R_{ff} = R_{ff} R_{ff}^H$, hence unitary eigenmatrix exists

Karhunen-Loeve transform

- Unitary transform with matrix

$$A = \Phi^H$$

where the columns of Φ are ordered according to decreasing eigenvalues.

- Transform coefficients are pairwise uncorrelated

$$R_{cc} = AR_{ff}A^H = \Phi^H R_{ff}\Phi = \Phi^H \Phi\Lambda = \Lambda$$

- Energy concentration property:
  - No other unitary transform packs as much energy into the first *J* coefficients, where *J* is arbitrary
  - Mean squared approximation error by choosing only first *J* coefficients is minimized.

Optimum energy concentration by KL transform

- To show optimum energy concentration property, consider the truncated coefficient vector

$$\vec{b} = I_J \vec{c}$$

where I_J contain ones on the first J diagonal positions, else zeros.

- Energy in first J coefficients for arbitrary transform A

$$E = Tr(R_{bb}) = Tr(I_J R_{cc} I_J) = Tr(I_J A R_{ff} A^H I_J) = \sum_{k=0}^{J-1} a_k^T R_{ff} a_k^*$$

where a_k^T is the k - th row of A.

- Lagrangian cost function to enforce unit-length basis vectors

$$L = E + \sum_{k=0}^{J-1} \lambda_k \left(1 - a_k^T a_k^*\right) = \sum_{k=0}^{J-1} a_k^T R_{ff} a_k^* + \sum_{k=0}^{J-1} \lambda_k \left(1 - a_k^T a_k^*\right)$$

- Differentiating L with respect to a_j yields neccessary condition

$$R_{ff} a_j^* = \lambda_j a_j^* \quad \text{for all } j < J$$

Basis images and eigenimages

- For any unitary transform, the inverse transform

$$\vec{f} = A^H \vec{c}$$

  can be interpreted in terms of the superposition of „basis images“ (columns of A^H) of size MN.
- If the transform is a KL transform, the basis images, which are the eigenvectors of the autocorrelation matrix R_{ff} , are called „eigenimages.“
- If energy concentration works well, only a limited number of eigenimages is needed to approximate a set of images with small error. These eigenimages form an optimal linear subspace of dimensionality J.

Computing eigenimages from a training set

- How to measure *MNxMN* autocorrelation matrix?
  - Use training set $\vec{\Gamma}_1, \vec{\Gamma}_2, \ldots, \vec{\Gamma}_L$
  - Define training set matrix $S = \left(\vec{\Gamma}_1, \vec{\Gamma}_2, \ldots, \vec{\Gamma}_L\right)$ and calculate
$$R_{\Gamma\Gamma} = \frac{1}{L}\sum_{l=1}^{L} \vec{\Gamma}_l \vec{\Gamma}_l^H = \frac{1}{L} SS^H$$
- **Problem 1:** Training set size should be $L >> MN$

  If $L < MN$, autocorrelation matrix R_{ff} is rank - deficient
- **Problem 2:** Finding eigenvectors of an *MNxMN* matrix.

- Can we find a small set of the most important eigenimages from a small training set $L << MN$

Sirovich and Kirby method

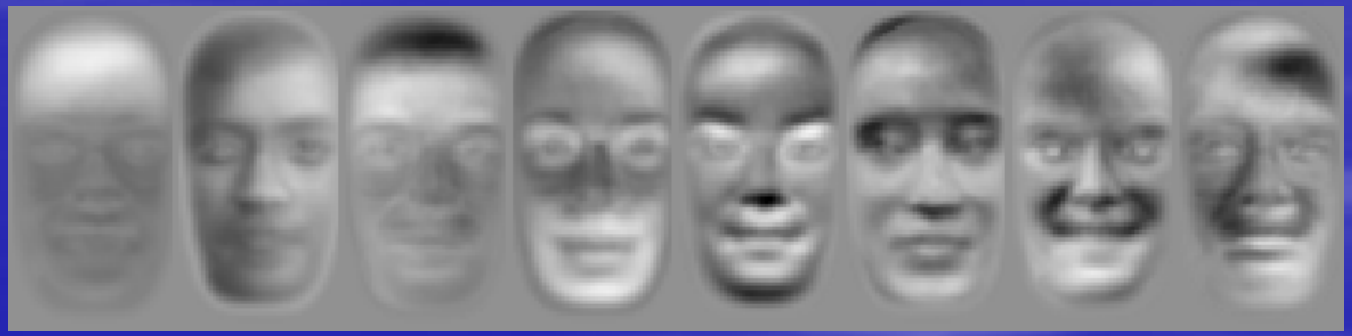
- Instead of eigenvectors of SS^H, consider the eigenvectors of S^HS, i.e.,

$$S^H S\vec{v}_i = \mu_i \vec{v}_i$$

- Premultiply both sides by S: $SS^H S\vec{v}_i = \mu_i S\vec{v}_i$
- By inspection, we find that $S\vec{v}_i$ are eigenvectors of SS^H
- For this gives rise to great computational savings, by
  - Computing the LxL matrix S^HS
  - Computing L eigenvectors $S\vec{v}_i$ of S^HS
  - Computing eigenimages corresponding to the $L_0 \leq L$ largest eigenvalues according as $S\vec{v}_i$

Example: eigenfaces

- The first 8 eigenfaces obtained from a training set of 500 frontal views of human faces.



- Can be used for face recognition by nearest neighbor search in 8-d „face space.“
- Can be used to generate faces by adjusting 8 coefficients.

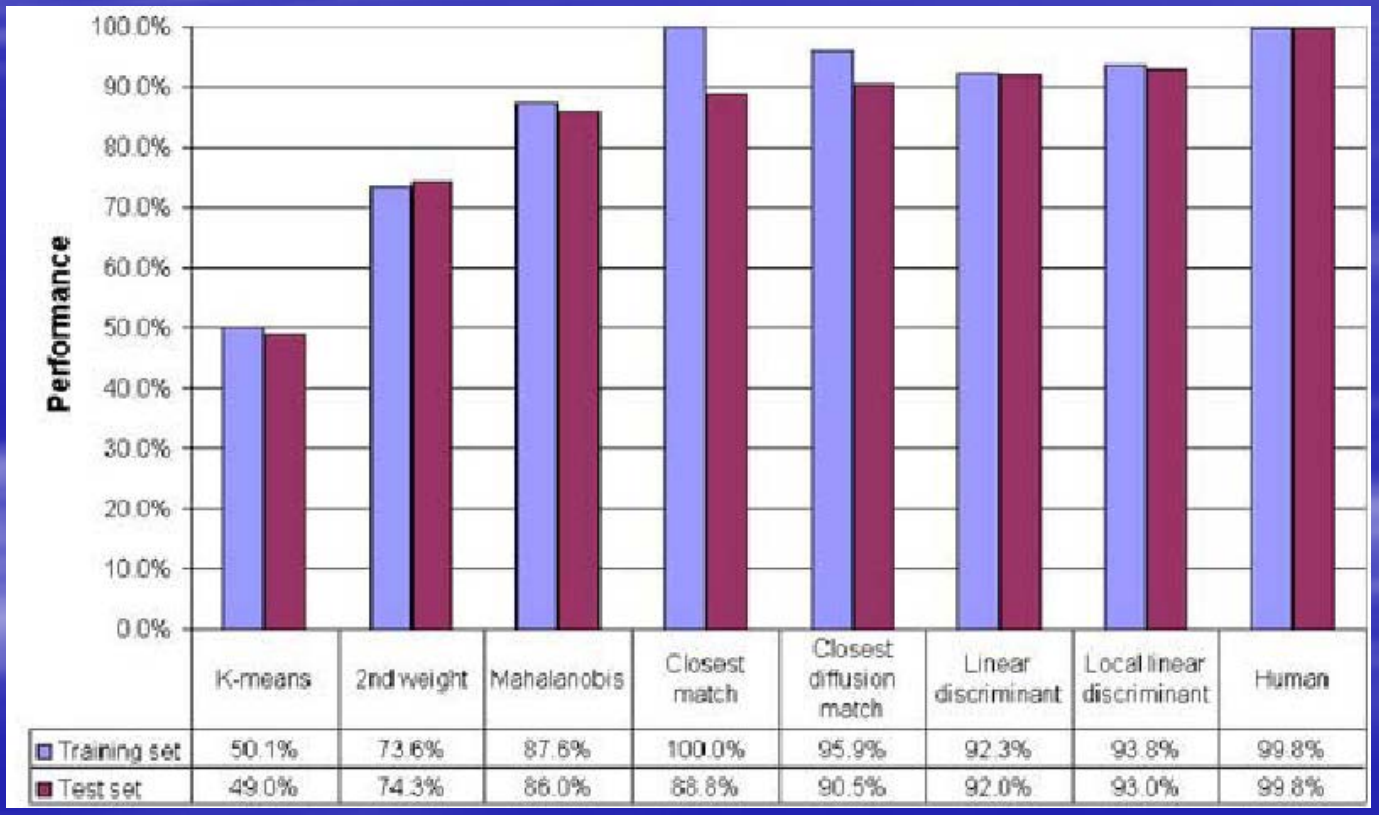
Gender recognition using eigenfaces

- Task: Male or female?

- Eigenimages from a data base of 20 and 20 female training images

Gender recognition using eigenfaces (cont.)

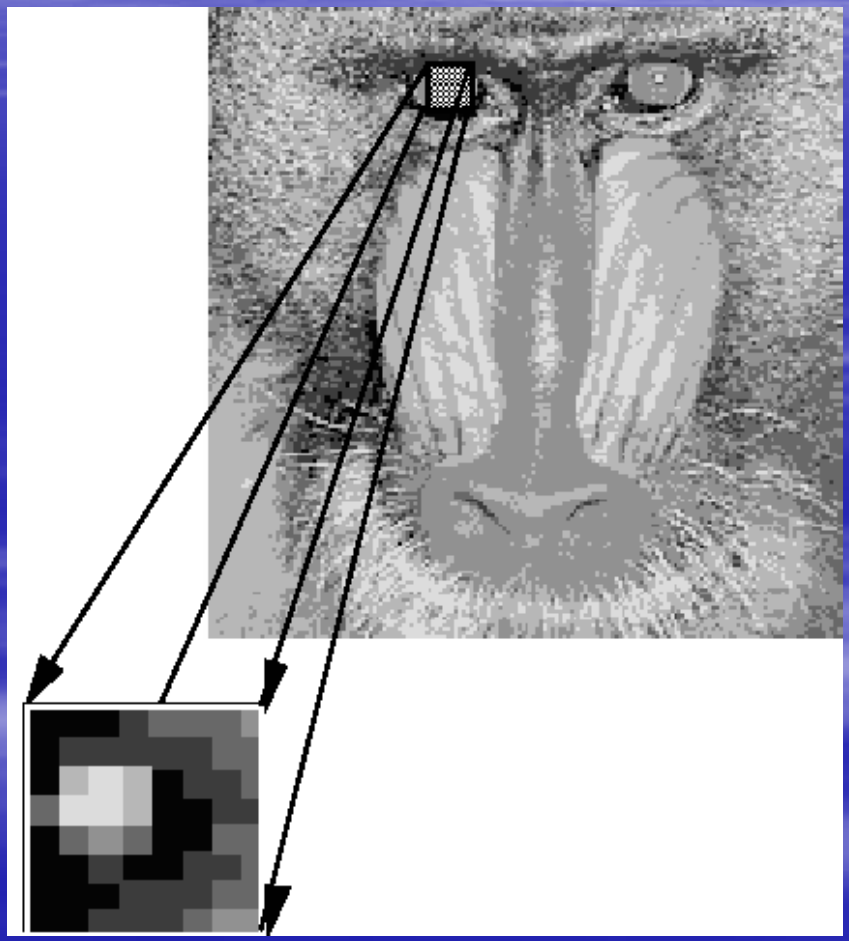
- Recognition accuracy using 8 eigenimages



	K-means	2nd weight	Mahalanobis	Closest match	Closest diffusion match	Linear discriminant	Local linear discriminant	Human
Training set	50.1%	73.6%	87.6%	100.0%	95.9%	92.3%	93.8%	99.8%
Test set	49.0%	74.3%	86.0%	68.8%	90.5%	92.0%	93.0%	99.8%

Block-wise image processing

- Subdivide image into small blocks
- Process each block independently from the others
- Typical blocksizes: 8x8, 16x16



Separable blockwise transforms

- Image block written as a square matrix $\mathbf{f}$

$$\mathbf{c} = \mathbf{A}^{\mathrm{T}}.\mathbf{f}.\mathbf{A}$$

(NxN coefficients)

- This can only be done, if transform is separable in *x* and *y*, i.e.,

$$\mathbf{A}^{\mathrm{T}} = \mathbf{A} \otimes \mathbf{A}$$

(N^2xN^2)

- Inverse transform

$$\mathbf{f} = \mathbf{A}^{*}.\mathbf{c}.\mathbf{A}^{\mathrm{H}}$$

Haar transform

- Haar transform matrix for sizes *N=2,4,8*

$$Hr_2 = \frac{1}{\sqrt{2}}\begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}$$

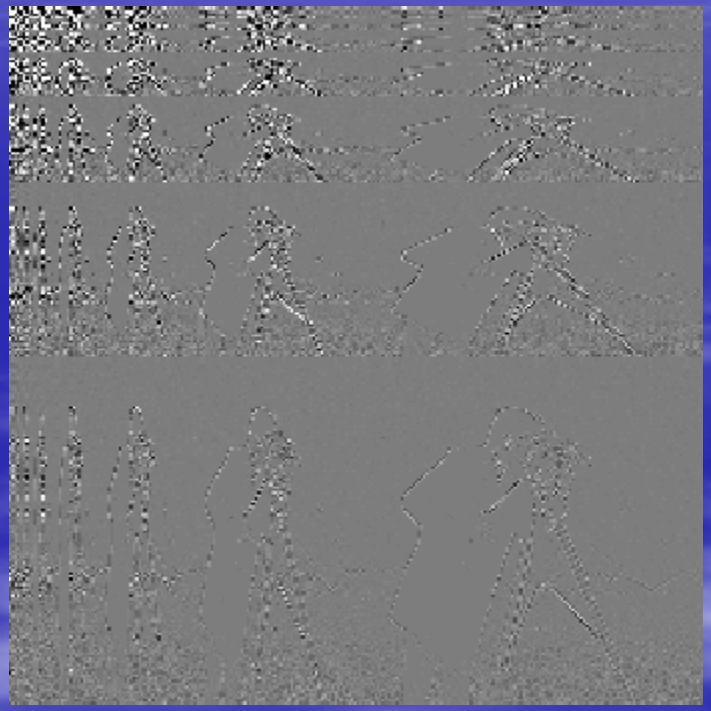
$$Hr_4 = \frac{1}{\sqrt{4}}\begin{pmatrix} 1 & 1 & \sqrt{2} & 0 \\ 1 & 1 & -\sqrt{2} & 0 \\ 1 & -1 & 0 & \sqrt{2} \\ 1 & -1 & 0 & -\sqrt{2} \end{pmatrix}$$

$$Hr_8 = \frac{1}{\sqrt{8}}\begin{pmatrix} 1 & 1 & \sqrt{2} & 0 & 2 & 0 & 0 & 0 \\ 1 & 1 & \sqrt{2} & 0 & -2 & 0 & 0 & 0 \\ 1 & 1 & -\sqrt{2} & 0 & 0 & 2 & 0 & 0 \\ 1 & 1 & -\sqrt{2} & 0 & 0 & -2 & 0 & 0 \\ 1 & -1 & 0 & \sqrt{2} & 0 & 0 & 2 & 0 \\ 1 & -1 & 0 & \sqrt{2} & 0 & 0 & -2 & 0 \\ 1 & -1 & 0 & -\sqrt{2} & 0 & 0 & 0 & 2 \\ 1 & -1 & 0 & -\sqrt{2} & 0 & 0 & 0 & -2 \end{pmatrix}$$

- Can be computed by taking sums and differences
- Fast algorithms by recursively applying $\mathbf{Hr_2}$

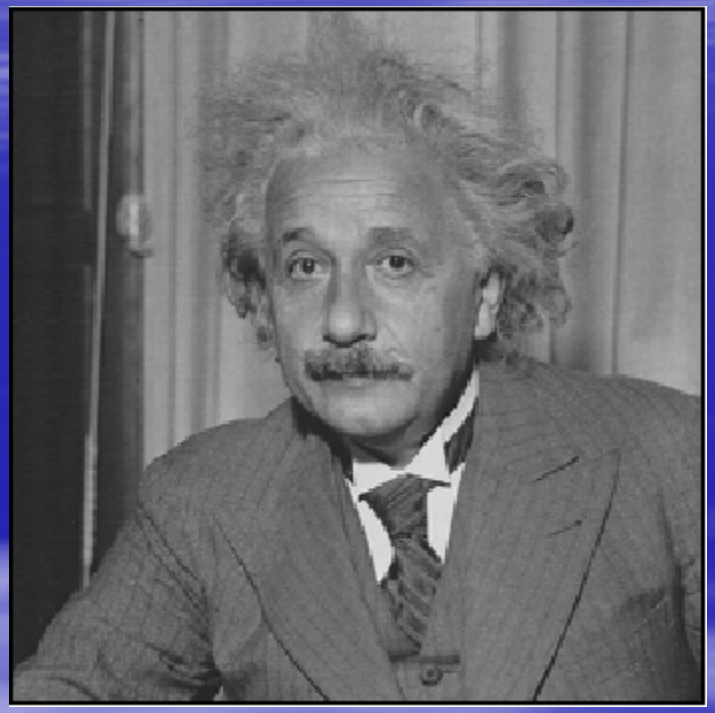
Haar transform example

Original *Cameraman*
256x256

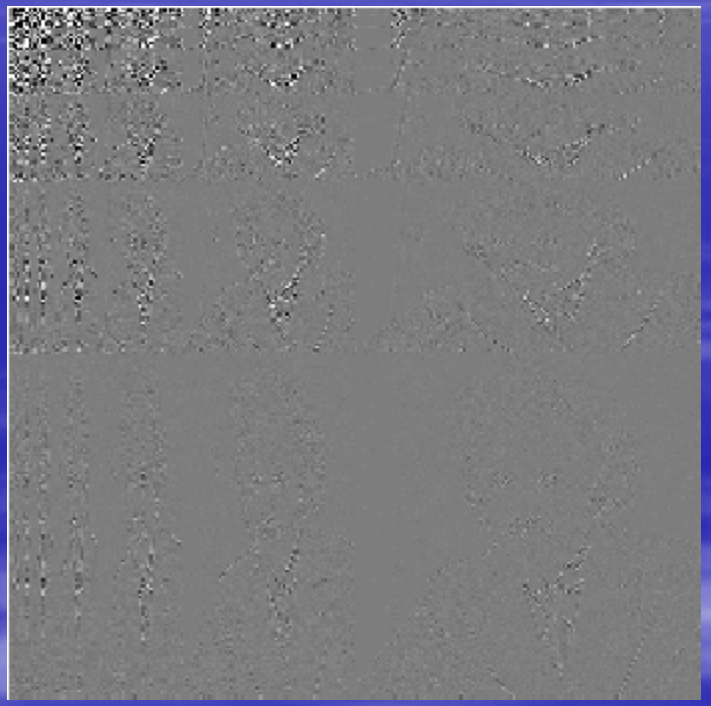


256x256 Haar transform
of *Cameraman*

Haar transform example



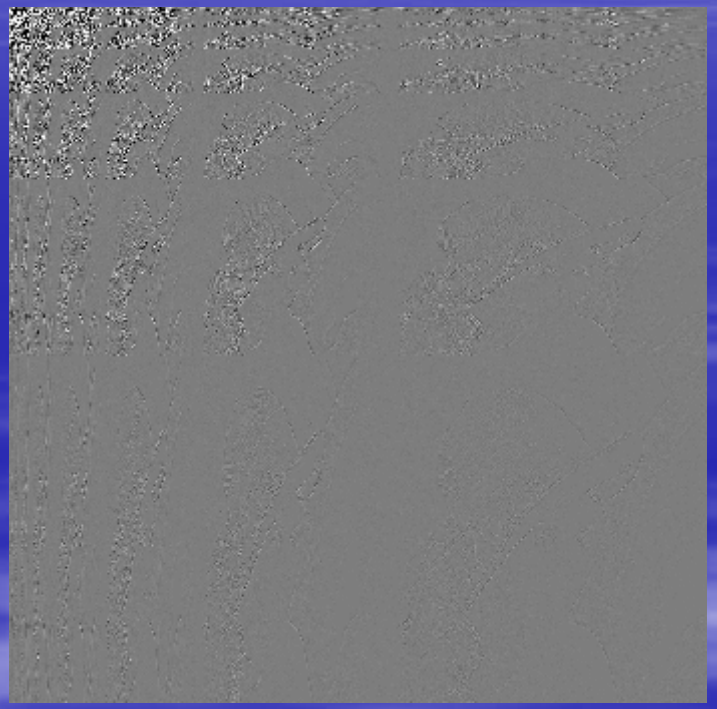
Original *Einstein*
256x256



256x256 Haar transform
of *Einstein*

Haar transform example

Original *Lena*
512x512



512x512 Haar transform
of *Lena*

Hadamard transform

- Transform matrices can be recursively generated

$$Hd_2 = Hr_2 = \frac{1}{\sqrt{2}}\begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}$$

$$Hd_4 = Hd_2 \otimes Hd_2$$

$$Hd_8 = Hd_4 \otimes Hd_2 = Hd_2 \otimes Hd_2 \otimes Hd_2$$

$$\dots$$

- Example

$$Hd_8 = \frac{1}{\sqrt{8}}\begin{pmatrix} 1 & 1 & 1 & 1 & 1 & 1 & 1 & 1 \\ 1 & -1 & 1 & -1 & 1 & -1 & 1 & -1 \\ 1 & 1 & -1 & -1 & 1 & 1 & -1 & -1 \\ 1 & -1 & -1 & 1 & 1 & -1 & -1 & 1 \\ 1 & 1 & 1 & 1 & -1 & -1 & -1 & -1 \\ 1 & -1 & 1 & -1 & -1 & 1 & -1 & 1 \\ 1 & 1 & -1 & -1 & -1 & -1 & 1 & 1 \\ 1 & -1 & -1 & 1 & -1 & 1 & 1 & -1 \end{pmatrix}$$

Note that Hadamard Coefficients need reordering to concentrate energy

Discrete Fourier transform

- For DFT of order *N*, define $W_N = \exp\left\{\frac{-j2\pi}{N}\right\}$
- Transform matrix

$$\mathrm{DTF}_N = \frac{1}{\sqrt{N}} \begin{pmatrix} W_N^0 & W_N^0 & W_N^0 & \cdots & W_N^0 \\ W_N^0 & W_N^1 & W_N^2 & \cdots & W_N^{N-1} \\ W_N^0 & W_N^2 & & \cdots & W_N^{2(N-1)} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ W_N^0 & W_N^{N-1} & W_N^{2(N-1)} & \cdots & W_N^{(N-1)(N-1)} \end{pmatrix}$$

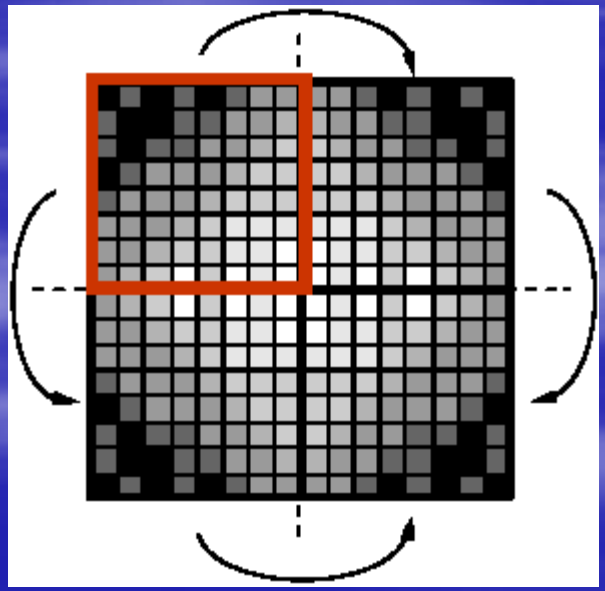
- Definition for general *N*, fast algorithms for $N=2^m$
- DFT coefficients are complex (even if input image is real)
- Inverse transform $\mathbf{DFT_N^{-1} = DTF^H = DTF^*}$

Discrete cosine transform

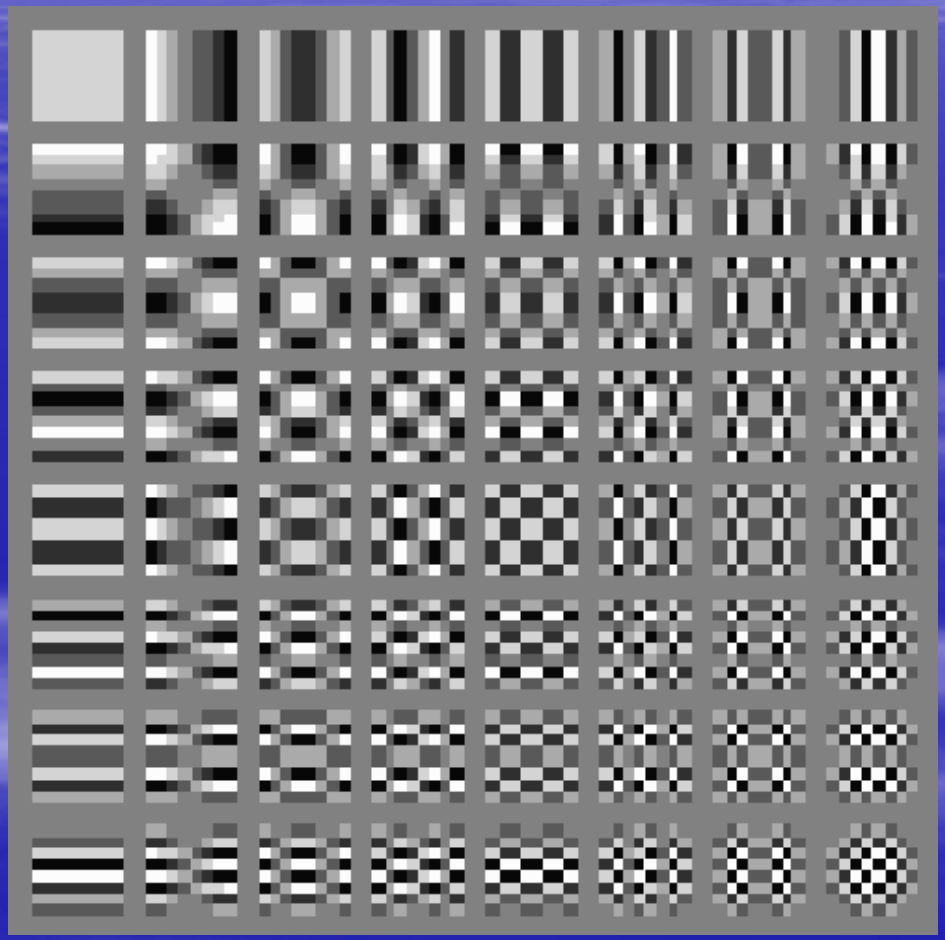
- Transform matrix

$$DCT_N = \left\{ a_{k,n} \middle| a_{k,n} = \begin{cases} \frac{1}{N} & 0 \le n \le N-1 \wedge k = 0 \\ \sqrt{\frac{2}{N}} \cos\left(\frac{\pi(2n+1)k}{2N} \right) & 0 \le n \le N-1 \wedge 1 \le k \le N-1 \end{cases} \right\}$$

- Can be interpreted as DFT of a mirror-extended image block (shown here for 2-d DCT)
- Transform is used in many coding standards (JPEG, MPEG)

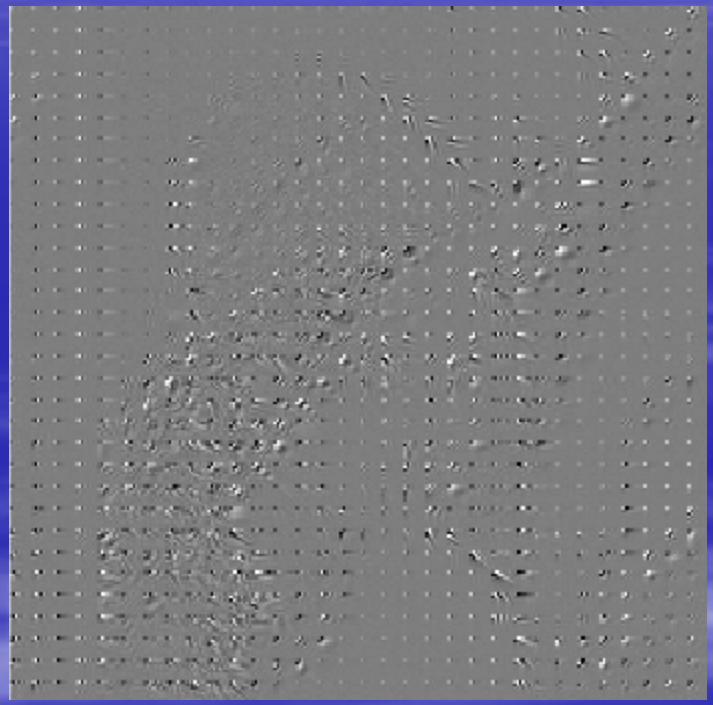


Basis images of an 8x8 DCT



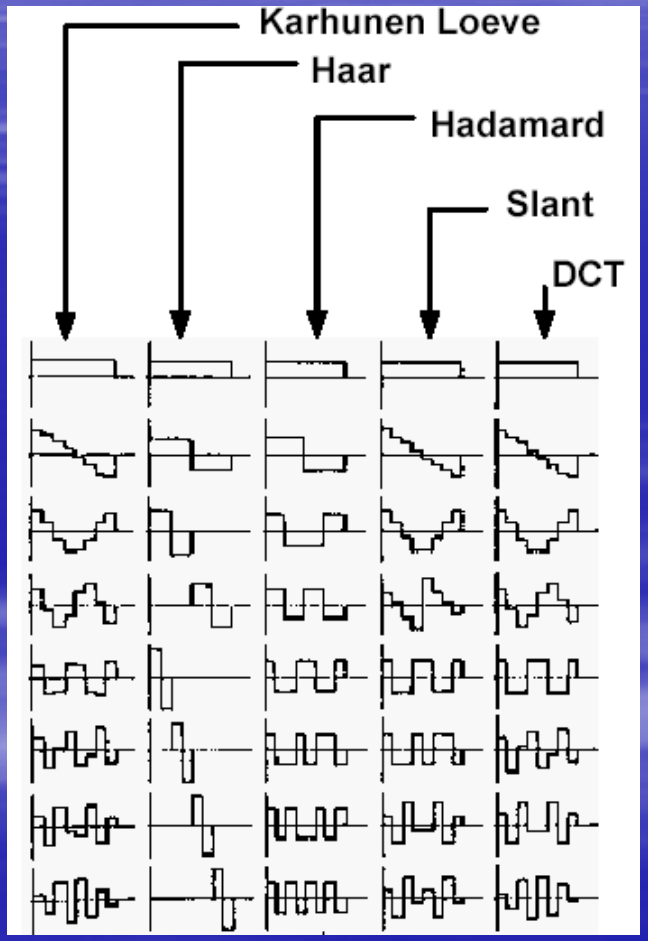
Blockwise DCT example

Original *Lena*
256x256



Blockwise 8x8 DCT
of *Lena*

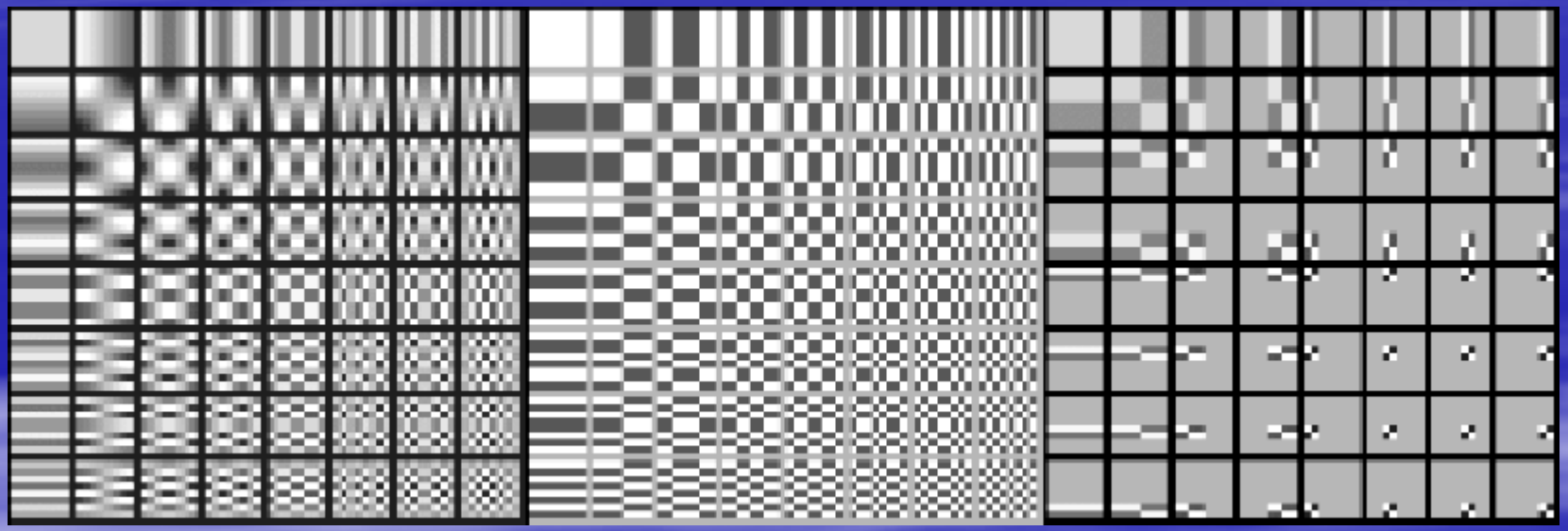
Comparison of block transforms



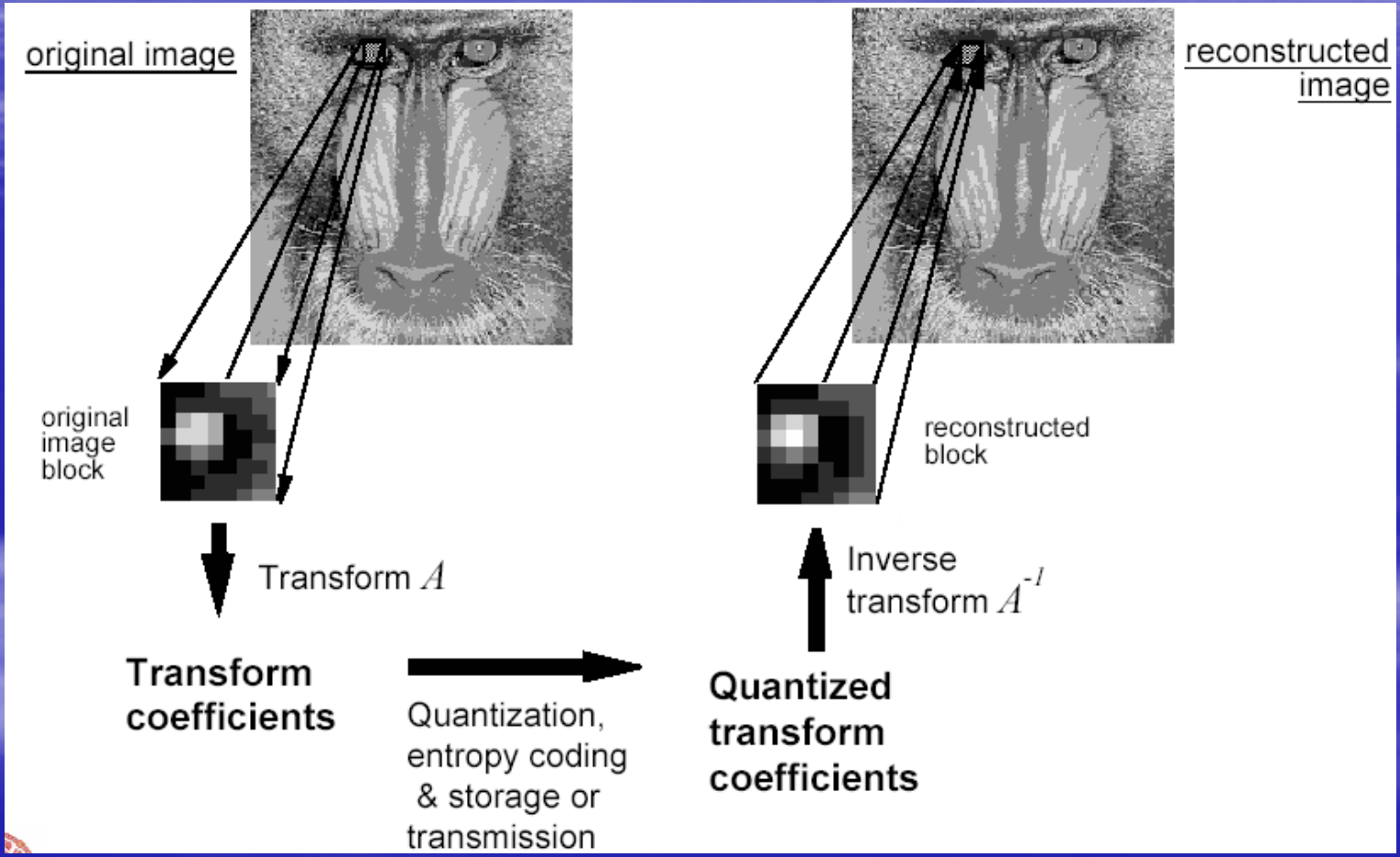
Comparison of 1-D basis functions for block size N=8

Comparison of block transforms (cont.)

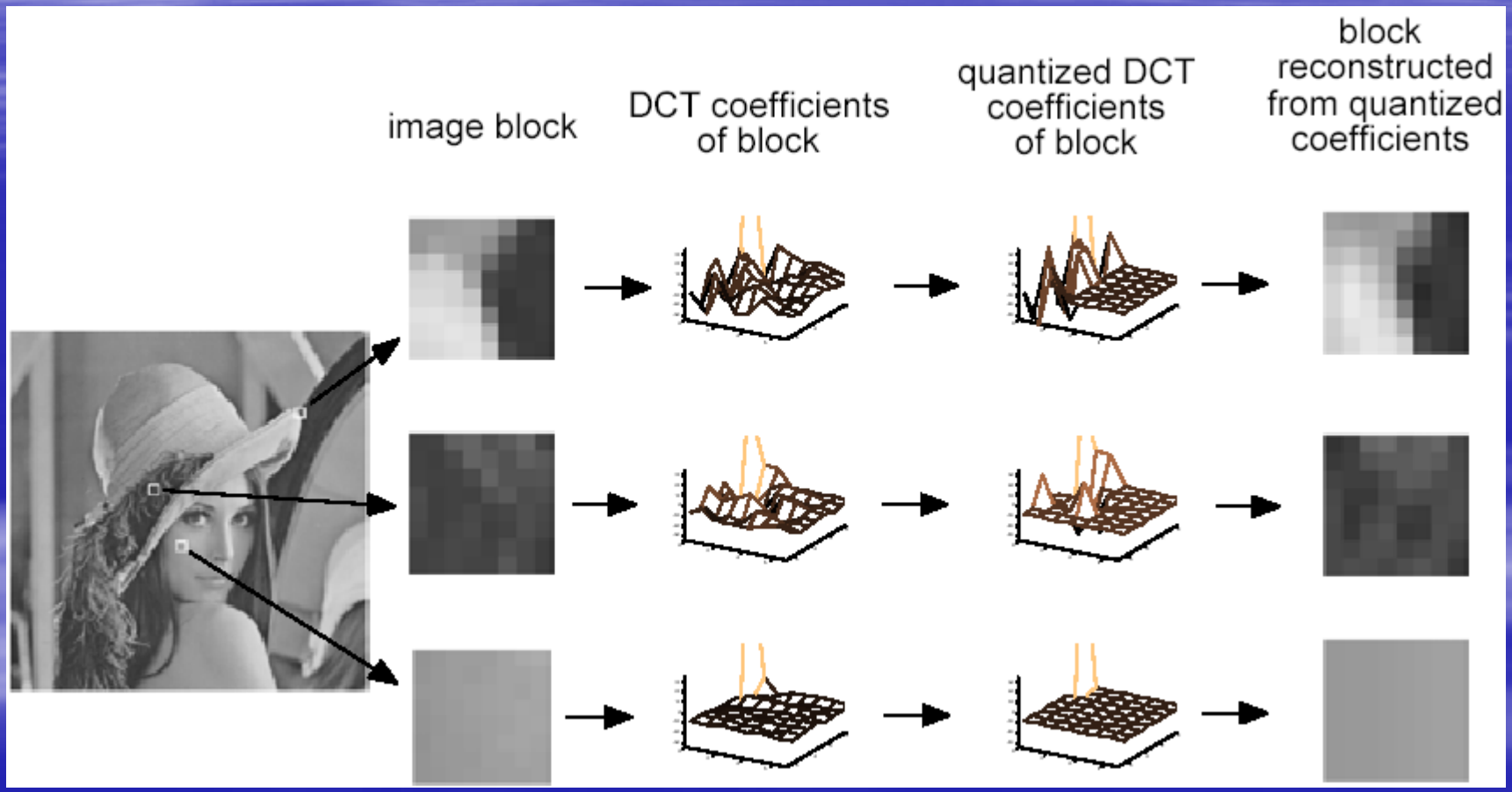
DCT Hadamard Haar



Transform Coding



Transform Coding (cont.)



DCT coding artifacts

DCT coding with increasingly coarse quantization, block size 8x8

quantizer stepsize
for AC coefficients: 25

quantizer stepsize
for AC coefficients: 100

quantizer stepsize
for AC coefficients: 200