Parsimonious Gaussian process models for the spectral-spatial classification of hyperspectral remote sensing images Seminar MIAT

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Nature of remote sensing images

A remote sensing image is a sampling of a spatial, spectral and temporel process







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Nature of remote sensing images

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Nature of remote sensing images

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Introduction

Hyperspectral Imagery 1/3





- Pixels are represented by random vector $\mathbf{x} \in \mathbb{R}^d$ with d large, associated to a random variable x that represents the class/label.
- Classification: predict the membership y of \mathbf{x} , $y = f(\mathbf{x})$.

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Hyperspectral Imagery 2/3

Instrument	Range (nm)	# Bands	Bandwidth (nm)	Spatial resolution (m)
AVIRIS	400-2500	224	10	20/1-4
HYDICE	400-2500	210	10	1-4
ROSIS-03	400-900	115	4	1
Hyspec	400-2500	427	3	1
HyMAP	400-2500	126	10-20	5
CASI	380-1050	288	2.4	1-2
HYPERION	400-2500	200	10	30

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Hyperspectral Imagery 3/3

Definition of more classes with finer resolution:



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Classification of hyperspectral ima

Image classification in high dimensional space

- High number of measurements but limited number of training samples.
- Curse of dimensionality: Statistical, geometrical and computational issues. Conventional method failed [Jimenez and Landgrebe, 1998].
- Kernel methods have shown great potential in many situations.
- Pixelwise classification not adapted [Fauvel et al., 2013].



Need to incorporate spatial information in the classification process: additional complexity.

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Kernel methods VS Parametric methods

- 1. Kernel methods [Camps-Valls and Bruzzone, 2009]:
 - Good abilities for classification,
 - Spatial information included through kernel function or additional features.

$$k_s(\mathbf{x}_i, \mathbf{x}_j) = \sum_{\substack{m \sim i \\ n \sim j}} k(\mathbf{x}_m, \mathbf{x}_n)$$

- 2. Parametric methods [Solberg et al., 1996]:
 - Markov Random Field: able to model spatial relationship between pixels,
 - Problem of the estimation of the spectral energy term.
- 3. Parametric kernel methods: probabilistic models in the kernel feature space.
 - Allow to get probability membership, with robust classifier
 - Allow to use the MRF modelization

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Kernel methods and MRF

- Maximum a posteriori: $\max_{Y}(Y|\mathbf{X})$
- When Y is MRF: $P(Y|\mathbf{X}) \propto \exp(-U(Y|\mathbf{X}))$ where $U(Y|\mathbf{X}) = \sum_{i=1}^{n} U(y_i|\mathbf{x}_i, \mathcal{N}_i)$ with

$$U(y_i|\mathbf{x}_i, \mathcal{N}_i) = \Omega(\mathbf{x}_i, y_i) + \rho \frac{\mathcal{E}(y_i, \mathcal{N}_i)}{\mathcal{E}(y_i, \mathcal{N}_i)}$$

• Spectral term: $-\log[p(\mathbf{x}_i|y_i)] \longleftarrow$

- SVM outputs [Farag et al., 2005, Tarabalka et al., 20/10, Moser and Serpico, 2013]
- Kernel-probabilistic model [Dundar and Landgrebe, 2004]

■ Spatial term ←

▶ Potts model: $\mathcal{E}(y_i, \mathcal{N}_i) = \sum_{j \in \mathcal{N}_i} [1 - \delta(y_i, y_j)]$

y_1	y_2	y_3
y_4	y_i	y_5
y_6	y_7	y_8

(Kernel) Gaussian mixture models

Quadratic decision rule in the input space

$$D_c(\mathbf{x}_i) = (\mathbf{x}_i - \boldsymbol{\mu}_c)^\top \boldsymbol{\Sigma}_c^{-1} (\mathbf{x}_i - \boldsymbol{\mu}_c) + \log(\det(\boldsymbol{\Sigma}_c)) - 2\ln(\pi_c)$$

Quadratic decision rule in the feature space [Dundar and Landgrebe, 2004]:

$$D_c(\phi(\mathbf{x}_i)) = \bar{\phi}_c(\mathbf{x}_i)^\top \mathbf{K}_c^{-1} \bar{\phi}_c(\mathbf{x}_i) + \log(\det(\mathbf{K}_c)) - 2\ln(\pi_c)$$

- Problem: K is badly conditioned (and non-invertible).
- Unlike SVM, there is no regularization for \mathbf{K}_c^{-1} and $\log(\det(\mathbf{K}_c))$ in the estimation process.
- So it needs to be included in the model.

(Kernel) Gaussian mixture models

Quadratic decision rule in the input space

$$D_c(\mathbf{x}_i) = (\mathbf{x}_i - \boldsymbol{\mu}_c)^\top \boldsymbol{\Sigma}_c^{-1} (\mathbf{x}_i - \boldsymbol{\mu}_c) + \log(\det(\boldsymbol{\Sigma}_c)) - 2\ln(\pi_c)$$

Quadratic decision rule in the feature space [Dundar and Landgrebe, 2004]:

$$D_c \left(\phi(\mathbf{x}_i)
ight) = ar{\phi}_c(\mathbf{x}_i)^\top \mathbf{K}_c^{-1} ar{\phi}_c(\mathbf{x}_i) + \log(\det(\mathbf{K}_c)) - 2\ln(\pi_c)$$

- Problem: **K** is badly conditioned (and non-invertible).
- Unlike SVM, there is no regularization for \mathbf{K}_c^{-1} and $\log(\det(\mathbf{K}_c))$ in the estimation process.
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Enforce parsimony in the model

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Gaussian process in the fea	ture space	

Kernel induced feature space



- Gaussian kernel: $k(\mathbf{x}_i, \mathbf{x}_j) = \exp\left(-\gamma \|\mathbf{x}_i \mathbf{x}_j\|_{\mathbb{R}^d}^2\right)$
- From Mercer theorem: $k(\mathbf{x}_i, \mathbf{x}_j) = \langle \phi(\mathbf{x}_i), \phi(\mathbf{x}_j) \rangle_{\mathcal{F}}$ which can be written

$$k(\mathbf{x}_i, \mathbf{x}_j) = \sum_{m=1}^{d_F} \lambda_m \mathbf{q}_m(\mathbf{x}_i) \mathbf{q}_m(\mathbf{x}_j)$$

where $d_{\mathcal{F}} = \dim(\mathcal{F})$.

$$\phi: \mathbf{x} \mapsto [\dots, \sqrt{\lambda_m} \mathbf{q}_m(\mathbf{x}), \dots], \ m = 1, 2, \dots, d_{\mathcal{F}}$$

 \blacksquare For the Gaussian kernel, $d_{\mathcal{F}}=+\infty$

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Gaussian process

- Let us assume that $\phi(\mathbf{x})$, conditionally on y = c, is a Gaussian process with mean μ_c and covariance function Σ_c .
- The projection of $\phi(\mathbf{x})$ on the eigenfunction \mathbf{q}_{cj} is noted $\phi(\mathbf{x})_j$:

$$\langle \phi(\mathbf{x}), \mathbf{q}_{cj} \rangle = \int_J \phi(\mathbf{x})(t) \mathbf{q}_{cj}(t) dt.$$

- The random vector $[\phi(\mathbf{x})_1, \ldots, \phi(\mathbf{x})_r] \in \mathbb{R}^r$ is, conditionally on y = c, a multivariate normal vector.
- Gaussian mixture model (Quadratic Discriminant) decision rules:

$$D_c(\phi(\mathbf{x}_i)) = \sum_{j=1}^r \left[\frac{\langle \phi(\mathbf{x}_i) - \boldsymbol{\mu}_c, \mathbf{q}_{cj} \rangle^2}{\lambda_{cj}} + \ln(\lambda_{cj}) \right] - 2\ln(\pi_c)$$

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Gaussian process

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- The random vector $[\phi(\mathbf{x})_1, \ldots, \phi(\mathbf{x})_r] \in \mathbb{R}^r$ is, conditionally on y = c, a multivariate normal vector.
- Gaussian mixture model (Quadratic Discriminant) decision rules: $r_c = \min(n_c, r)$

$$D_{c}(\phi(\mathbf{x}_{i})) = \sum_{j=1}^{r_{c}} \left[\frac{\langle \phi(\mathbf{x}_{i}) - \boldsymbol{\mu}_{c}, \mathbf{q}_{cj} \rangle^{2}}{\lambda_{cj}} + \ln(\lambda_{cj}) \right] - 2\ln(\pi_{c}) \\ + \sum_{j=r_{c}+1}^{r} \left[\frac{\langle \phi(\mathbf{x}_{i}) - \boldsymbol{\mu}_{c}, \mathbf{q}_{cj} \rangle^{2}}{\lambda_{cj}} + \ln(\lambda_{cj}) \right]$$

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Definitions

Definition (Parsimonious Gaussian process with common noise)

 $p\mathcal{GP}$ is a Gaussian process $\phi(\mathbf{x})$ for which, conditionally to y = c, the eigen-decomposition of its covariance operator Σ_c is such that

A1. It exists a dimension $r < +\infty$ such that $\lambda_{cj} = 0$ for $j \ge r$ and for all $c = 1, \dots, C$.

A2. It exists a dimension $p_c < \min(r, n_c)$ such that $\lambda_{cj} = \lambda$ for $p_c < j < r$ and for all $c = 1, \ldots, C$.

Definition (Parsimonious Gaussian process with class specific noise)

- A3. It exists a dimension $r_c < r$ such that $\lambda_{cj} = 0$ for all $j > r_c$ and for all $c = 1, \ldots, C$. When $r = +\infty$, it is assumed that $r_c = n_c 1$.
- A4. It exists a dimension $p_c < r_c$ such that $\lambda_{cj} = \lambda_c$ for $j > p_c$ and $j \le r_c$, and for all $c = 1, \ldots, C$.
 - A1 and A3 are motivated by the quick decay of the eigenvalues of Gaussian kernels.

• A2 and A4 express that the data of each class lives in a specific subspace of size p_c .

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$p\mathcal{GP}$ models: List of sub-models

Model	Variance inside \mathcal{F}_c	\mathbf{q}_{cj}	p_c
	Variance outside \mathcal{F}_c : Common		
$p{\cal GP}_0$	Free	Free	Free
$p \mathcal{GP}_1$	Free	Free	Common
$p{\cal GP}_2$	Common within groups	Free	Free
$p{\cal GP}_3$	Common within groups	Free	Common
$p{\cal GP}_4$	Common between groups	Free	Common
$p{\cal GP}_5$	Common within and between groups	Free	Free
$p{\cal GP}_6$	Common within and between groups	Free	Common
	Variance outside \mathcal{F}_c : Free		
$np {\cal GP}_0$	Free	Free	Free
$np \mathcal{GP}_1$	Free	Free	Common
$np {\cal GP}_2$	Common within groups	Free	Free
$np {\cal GP}_3$	Common within groups	Free	Common
$np {\cal GP}_4$	Common between groups	Free	Common

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Figure: Visual illustration of model $np\mathcal{GP}_1$. Dimension of \mathcal{F}_c is common to both classes, they have specific variance inside \mathcal{F}_c and they have specific noise level.

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Decision rules for $p\mathcal{GP}_0$

Proposition

For $p\mathcal{GP}_0$, the decision rule can be written:

$$D_{c}(\phi(\mathbf{x}_{i})) = \sum_{j=1}^{p_{c}} \frac{\lambda - \lambda_{cj}}{\lambda_{cj}\lambda} \langle \phi(\mathbf{x}_{i}) - \boldsymbol{\mu}_{c}, \mathbf{q}_{cj} \rangle^{2} - 2\ln(\pi_{c}) + \frac{\|\phi(\mathbf{x}) - \boldsymbol{\mu}_{c}\|^{2}}{\lambda} + \sum_{j=1}^{p_{c}} \ln(\lambda_{cj}) + (p_{M} - p_{c})\ln(\lambda) + \gamma$$

where γ is a constant term that does not depend on the index c of the class.

- Proofs are given in [Bouveyron et al., 2014].
- \blacksquare Decompose the sum: $\sum_{j=1}^{p_c}\lambda_{cj}+\sum_{j=p_c+1}^r\lambda$
- \blacksquare Use the property: $\sum_{j=1}^r \langle \phi(\mathbf{x}) \boldsymbol{\mu}_c, \mathbf{q}_{cj} \rangle^2 = \|\phi(\mathbf{x}) \boldsymbol{\mu}_c\|^2$

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Estimation of the parameters

• Centered Gaussian kernel function according to class *c*:

$$\bar{k}_{c}(\mathbf{x}_{i},\mathbf{x}_{j}) = k(\mathbf{x}_{i},\mathbf{x}_{j}) + \frac{1}{n_{c}^{2}} \sum_{\substack{l,l'=1\\y_{l},y_{l}'=c}}^{n_{c}} k(\mathbf{x}_{l},\mathbf{x}_{l'}) - \frac{1}{n_{c}} \sum_{\substack{l=1\\y_{l}=c}}^{n_{c}} \left(k(\mathbf{x}_{i},\mathbf{x}_{l}) + k(\mathbf{x}_{j},\mathbf{x}_{l})\right).$$

and
$$\overline{\mathbf{K}}_c$$
 of size $n_c \times n_c$: $(\overline{\mathbf{K}}_c)_{l,l'} = \frac{k_c(\mathbf{x}_l, \mathbf{x}_{l'})}{n_c}$.

• $\hat{\lambda}_{cj}$ is the j^{th} largest eigenvalue of $\overline{\mathbf{K}}_{c}$, and $\boldsymbol{\beta}_{cj}$ is its associated normalized eigenvector.

$$\hat{\lambda} = \frac{1}{\sum_{c=1}^{C} \hat{\pi}_c (r_c - \hat{p}_c)} \sum_{c=1}^{C} \hat{\pi} \left(\operatorname{trace}(\overline{\mathbf{K}}_c) - \sum_{j=1}^{\hat{p}_c} \hat{\lambda}_{cj} \right).$$

$$\hat{\pi}_c = n_c/n.$$

• \hat{p}_c : percentage of cumulative variance.

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Computable decision rule

Proposition

The decision rule can be computed as:

$$D_c(\phi(\mathbf{x}_i)) = \frac{1}{n_c} \sum_{j=1}^{\hat{p}_c} \frac{\hat{\lambda} - \hat{\lambda}_{cj}}{\hat{\lambda}_{cj}^2 \hat{\lambda}} \left(\sum_{\substack{l=1\\y_l=c}}^{n_c} \beta_{cjl} \bar{k}_c(\mathbf{x}_i, \mathbf{x}_l) \right)^2 + \frac{\bar{k}_c(\mathbf{x}_i, \mathbf{x}_i)}{\hat{\lambda}} + \sum_{j=1}^{\hat{p}_c} \ln(\hat{\lambda}_{cj}) + (\hat{p}_M - \hat{p}_c) \ln(\hat{\lambda}) - 2 \ln(\hat{\pi}_c)$$

Proofs are given in [Bouveyron et al., 2014].

- Use of the property that the eigenfunction of the covariance function is a linear combination of $\phi({\bf x}_i)-{\pmb \mu}_c$

$$\langle \phi(\mathbf{x}_i) - \boldsymbol{\mu}_c, \phi(\mathbf{x}_j) - \boldsymbol{\mu}_c \rangle = \bar{k}_c(\mathbf{x}_i, \mathbf{x}_j)$$

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Numerical considerations

- The proposed model allow a *safe* computation of \mathbf{K}_c^{-1} and $\log (\det(\mathbf{K}_c))$ that appears in the kernel quadratic decision rule.
- Only the p_c first eigenvector/eigenvalue are used
- Eigenvectors corresponding to small eigenvalues are not used
- If p_c s are not too large, $\log(\hat{\lambda})$ is stable.

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Proof: \mathbf{K}_c is *pdf* so it can be decomposed into $\mathbf{Q}_c \mathbf{\Lambda}_c \mathbf{Q}_c^{\top} = \sum_{j=1}^r \lambda_{cj} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top}$

$$\begin{aligned} \mathbf{K}_{c}^{-1} &= \mathbf{Q}_{c} \mathbf{\Lambda}_{c}^{-1} \mathbf{Q}_{c}^{\top} = \sum_{j=1}^{r} \lambda_{cj}^{-1} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} = \sum_{j=1}^{p_{c}} \lambda_{cj}^{-1} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} + \lambda^{-1} \sum_{j=p_{c}+1}^{r} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} \\ &= \sum_{j=1}^{p_{c}} \lambda_{cj}^{-1} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} + \lambda^{-1} \left(\mathbf{I}_{n_{c}} - \sum_{j=1}^{p_{c}} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} \right) = \sum_{j=1}^{p_{c}} \frac{\lambda - \lambda_{cj}}{\lambda \lambda_{cj}} \mathbf{q}_{cj} \mathbf{q}_{cj}^{\top} + \lambda^{-1} \mathbf{I}_{n_{c}} \end{aligned}$$

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Numerical considerations

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$$\mathbf{K}_{c}^{-1} = \mathbf{Q}_{c}\mathbf{\Lambda}_{c}^{-1}\mathbf{Q}_{c}^{\top} = \sum_{j=1}^{r} \lambda_{cj}^{-1}\mathbf{q}_{cj}\mathbf{q}_{cj}^{\top} = \sum_{j=1}^{p_{c}} \lambda_{cj}^{-1}\mathbf{q}_{cj}\mathbf{q}_{cj}^{\top} + \lambda^{-1}\sum_{j=p_{c}+1}^{r} \mathbf{q}_{cj}\mathbf{q}_{cj}^{\top}$$
$$= \sum_{j=1}^{p_{c}} \lambda_{cj}^{-1}\mathbf{q}_{cj}\mathbf{q}_{cj}^{\top} + \lambda^{-1}\left(\mathbf{I}_{n_{c}} - \sum_{j=1}^{p_{c}} \mathbf{q}_{cj}\mathbf{q}_{cj}^{\top}\right) = \sum_{j=1}^{p_{c}} \frac{\lambda - \lambda_{cj}}{\lambda\lambda_{cj}}\mathbf{q}_{cj}\mathbf{q}_{cj}^{\top} + \lambda^{-1}\mathbf{I}_{n_{c}}$$

$$\log\left(\det(\mathbf{K}_c)\right) = \sum_{j=1}^{p_c} \log(\lambda_{cj}) + (r - p_c) \log(\lambda)$$

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Existing models

- [Dundar and Landgrebe, 2004] Equal covariance matrix assumption and ridge regularization. Complexity: $\mathcal{O}(n^3)$. Similar to $p\mathcal{GP}_4$ with equal eigenvectors.
- [Pekalska and Haasdonk, 2009] Ridge regularization, per class. Complexity: $\mathcal{O}(n_c^3)$.
- [Xu et al., 2009] The last $n_c - p - 1$ eigenvalues are equal to λ_{cp} . Complexity: $\mathcal{O}(n_c^3)$. Similar to $p\mathcal{GP}_1$.

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Existing models

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- [Pekalska and Haasdonk, 2009]
 Ridge regularization, per class. Complexity: O(n_c³).
- [Xu et al., 2009]

The last $n_c - p - 1$ eigenvalues are equal to λ_{cp} . Complexity: $\mathcal{O}(n_c^3)$. Similar to $p\mathcal{GP}_1$.



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Data sets

- University of Pavia: 103 spectral bands, 9 classes and 42,776 referenced pixels.
- Kennedy Space Center: 224 spectral bands, 13 classes and 4,561 referenced pixels.
- Heves: 252 spectral bands, 16 classes and 360,953 pixels.





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Protocol

- [Fauvel et al., 2015]
- **5**0 training pixels for each class have been randomly selected from the samples.
- The remaining set of pixels has been used for validation to compute the correct classification rate.
- Repeated 20 times.
- Variables have been scaled between 0 and 1.
- Competitive methods
 - SVM
 - RF
 - Kernel-DA (M. Dundar and D. A. Landgrebe, 2004)
- Hyperparameters learn by 5-CV.

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Classification accuracy

	Kappa coefficient		Process	ing time	e (s)	
	University	KSC	Heves	University	KSC	Heves
$p{\cal GP}_0$	0.768	0.920	0.664	18	31	148
$p{\cal GP}_1$	0.793	0.922	0.671	18	33	151
$p{\cal GP}_2$	0.617	0.844	0.588	18	31	148
$p{\cal GP}_3$	0.603	0.842	0.594	19	33	152
$p{\cal GP}_4$	0.661	0.870	0.595	19	34	152
$p{\cal GP}_5$	0.567	0.820	0.582	18	32	148
$p{\cal GP}_6$	0.610	0.845	0.583	19	34	152
$np {\cal GP}_0$	0.730	0.911	0.640	17	31	148
$np \mathcal{GP}_1$	0.792	0.921	0.677	18	33	151
$np {\cal GP}_2$	0.599	0.838	0.573	18	31	148
$np {\cal GP}_3$	0.578	0.817	0.585	19	33	152
$np \mathcal{GP}_4$	0.578	0.817	0.585	19	33	152
KDC	0.786	0.924	0.666	98	253	695
RF	0.646	0.853	0.585	3	3	18
SVM	0.799	0.928	0.658	10	28	171

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$p\mathcal{GP}\mathsf{MRF}$







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- Family of parsimonious Gaussian process models.
- Good performances wrt SVM and KDA
- Faster computation than previous KDA.
- $(n)p\mathcal{GP}_1$ perform the best.
- MRF extension.
- https://github.com/mfauvel/PGPDA
- Extension:
 - Non numerical data
 - Binary data
 - Unsupervised learning

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