Dictionary Learning Applications in Control Theory

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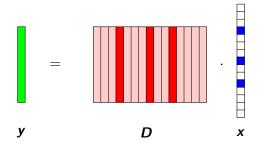
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Sparse Representation (SR)



Orthogonal Matching Pursuit (OMP)

Algorithm 1: OMP^a

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1 Arguments: \textbf{\textit{D}}, \textbf{\textit{y}}, \textbf{\textit{s}}

2 Initialize: \textbf{\textit{r}} = \textbf{\textit{y}}, \mathcal{I} = \emptyset

3 for k = 1: \textbf{\textit{s}} do

4 | Compute correlations with residual: \textbf{\textit{z}} = \textbf{\textit{D}}^T \textbf{\textit{r}}

5 | Select new column: i = \arg\max_{j}|z_j|

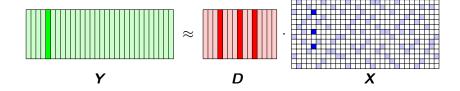
6 | Increase support: \mathcal{I} \leftarrow \mathcal{I} \cup \{i\}

7 | Compute new solution: \textbf{\textit{x}} = \mathsf{LS}(\textbf{\textit{D}}, \textbf{\textit{y}}, \mathcal{I})

8 | Update residual: \textbf{\textit{r}} = \textbf{\textit{y}} - \textbf{\textit{D}}_{\mathcal{I}} \textbf{\textit{x}}_{\mathcal{I}}
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^aPati, Rezaiifar, and Krishnaprasad 1993.

Dictionary Learning (DL)



The Dictionary Learning (DL) Problem

Given a data set $Y \in \mathbb{R}^{p \times m}$ and a sparsity level s, minimize the bivariate function

minimize
$$\| \boldsymbol{Y} - \boldsymbol{D} \boldsymbol{X} \|_F^2$$

subject to $\| \boldsymbol{d}_j \|_2 = 1, \ 1 \le j \le n$ (1) $\| \boldsymbol{x}_i \|_0 \le s, \ 1 \le i \le m,$

where $D \in \mathbb{R}^{p \times n}$ is the dictionary (whose columns are called atoms) and $X \in \mathbb{R}^{n \times m}$ the sparse representations matrix.

Approach

Algorithm 2: Dictionary learning - general structure

- 1 Arguments: signal matrix \boldsymbol{Y} , target sparsity s
- 2 Initialize: dictionary **D** (with normalized atoms)
- 3 for k = 1, 2, ... do
- 4 With fixed **D**, compute sparse representations **X**
- With fixed $m{X}$, update atoms $m{d}_j$, j=1:n

DL Algorithms

K-SVD¹ solves the optimization problem in sequence

$$\min_{\boldsymbol{d}_{j},\boldsymbol{X}_{j},\mathcal{I}_{j}} \left\| \left(\boldsymbol{Y}_{\mathcal{I}_{j}} - \sum_{\ell \neq j} \boldsymbol{d}_{\ell} \boldsymbol{X}_{\ell,\mathcal{I}_{\ell}} \right) - \boldsymbol{d}_{j} \boldsymbol{X}_{j,\mathcal{I}_{j}} \right\|_{F}^{2}$$
(2)

where all atoms excepting d_i are fixed.

This is seen as a rank-1 approximation and the solution is given by the singular vectors corresponding to the largest singular value.

$$\mathbf{d}_{j} = \mathbf{u}_{1}, \quad \mathbf{X}_{j,\mathcal{I}_{j}} = \sigma_{1}\mathbf{v}_{1}. \tag{3}$$



¹Aharon, Elad, and Bruckstein 2006.

LC-KSVD

minimize
$$\| \boldsymbol{Y} - \boldsymbol{D} \boldsymbol{X} \|_F^2 + \alpha \| \boldsymbol{Q} - \boldsymbol{A} \boldsymbol{X} \|_F^2 + \beta \| \boldsymbol{H} - \boldsymbol{W} \boldsymbol{X} \|_F^2$$
 subject to $\| \boldsymbol{d}_j \|_2 = 1, \ 1 \le j \le n$ (4) $\| \boldsymbol{x}_i \|_0 \le s, \ 1 \le i \le m,$

- dictionary atoms evenly split among classes
- q_i has non-zero entries where y_i and d_i share the same label.
- ullet linear transformation $oldsymbol{A}$ encourages discrimination in $oldsymbol{X}$
- $h_i = e_j$ where j is the class label of y_i
- W represents the learned classifier parameters

Fault Detection and Isolation in Water Networks

FDI via DL

Water networks pose some interesting issues:

- large scale, distributed network with few sensors
- user demand unknown or imprecise
- pressure dynamics nonlinear (analytic solutions impractical)

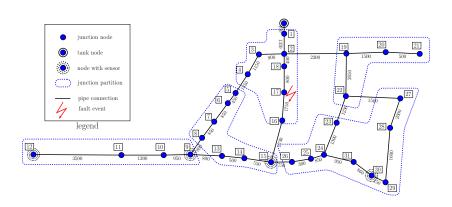
The DL approach for FDI:

a residual signal compares expected and measured pressures

$$r_i(t) = p_i(f_i(t), f_j(t), t) - \bar{p}_i, \forall i, j$$
 (5)

- to each fault is assigned a class and DL provides the atoms which discriminate between them
- each residual is sparsely described by atoms and thus, FDI is achieved iff the classification is unambiguous

Hanoi



Sensor Placement

Let $R \in \mathbb{R}^{n \times mn}$ be measured pressure residuals in all n network nodes. For each node we simulate m different faults.

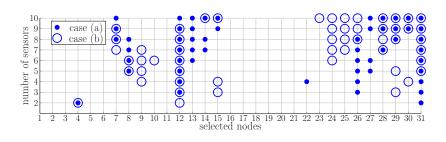
Given s < n available sensors, apply OMP on each column r

minimize
$$\| \mathbf{r} - \mathbf{I}_n \mathbf{x} \|_2^2$$
 subject to $\| \mathbf{x} \|_0 \le s$, (6)

resulting in matrix \boldsymbol{X} with s-sparse columns approximating \boldsymbol{R} .

Placement Strategies

- (a) select the s most common used atoms
- (b) from each m-block select most frequent s atoms; of the $n \cdot s$ atoms, select again the first s.



Learning

Algorithm 3: Placement and FDI learning^a

- 1 Inputs: training residuals $\mathbf{R} \in \mathbb{R}^{n \times nm}$
- 2 parameters s, α , β
- 3 Result: dictionary D, classifier W, sensor nodes \mathcal{I}_s
- 4 Select s sensor nodes \mathcal{I}_s based on matrix R using (a) or (b)
- 5 Let $extbf{\textit{R}}_{ extsf{\textit{I}}_{ extsf{\textit{s}}}}$ be the restriction of $extbf{\textit{R}}$ to the rows in $extsf{\textit{I}}_{ extsf{\textit{s}}}$
- 6 Use $R_{\mathcal{I}_s}$, α and β to learn D and W from (4)

^aIrofti and Stoican 2017.

Fault Detection

Algorithm 4: Fault detection and isolation

```
1 Inputs: testing residuals R \in \mathbb{R}^{s \times mn}

2 dictionary D, classifier W

3 Result: prediction P \in \mathbb{N}^{mn}

4 for k = 1 to mn do

5 Use OMP to obtain x_k using r_k and D

6 Label: L_k = Wx_k

7 Classify: p_k = \arg\max_c L_k
```

Position c of the largest entry from L_k is the predicted class.

Today

Improved sensor placement. Iteratively choose s rows from R solving at each step

$$i = \underset{k}{\operatorname{arg\,min}} \quad \|\operatorname{proj}_{\boldsymbol{R}_{\mathcal{I}}}\boldsymbol{r}_{k}\|_{2}^{2} + \lambda \frac{1}{\|\boldsymbol{\delta}_{k,\mathcal{I}}\|_{1}}, \boldsymbol{r}_{k} \in \boldsymbol{R}_{\mathcal{I}^{c}}, \tag{7}$$

where \mathcal{I} is the set of currently selected rows and $\delta_{k,\mathcal{I}}$ vector elements are the distances from node k to the nodes in \mathcal{I} .

Graph aware DL. Adding graph regularization²

$$\|\mathbf{Y} - \mathbf{D}\mathbf{X}\|_{F}^{2} + \alpha \|\mathbf{Q} - \mathbf{A}\mathbf{X}\|_{F}^{2} + \beta \|\mathbf{H} - \mathbf{W}\mathbf{X}\|_{F}^{2} + \gamma \text{Tr}(\mathbf{D}^{T}\mathbf{L}\mathbf{D}) + \lambda \text{Tr}(\mathbf{X}\mathbf{L}_{c}\mathbf{X}^{T}) + \mu \|\mathbf{L}\|_{F}^{2},$$
(8)

where \boldsymbol{L} is the graph Laplacian.



²Yankelevsky and Elad 2016.

Zonotopic Area Coverage

Zonotopic sets

Area packing, mRPI (over)approximation and other related notions may be described via unions of zonotopic sets:

$$\min_{Z_k} \operatorname{vol}(S) - \operatorname{vol}\left(\bigcup_k Z_k\right),$$
subject to $Z_k \subseteq S$.

Zonotopes, given in generator representation³

$$Z_k = \mathcal{Z}(c_k, G_k) = \{c_k + G_k \xi : ||\xi||_{\infty} \le 1\}$$
 (10)

are easy to handle for:

- Minkowski sum: $\mathcal{Z}(G_1, c_1) \oplus \mathcal{Z}(G_2, c_2) = \mathcal{Z}(\begin{bmatrix} G_1 & G_2 \end{bmatrix}, c_1 + c_2)$
- linear mappings: $RZ(G_1, c_1) = Z(RG_1, Rc_1)$



³Fukuda 2004.

Formulation

Each zonotope is parameterized after its center and a scaling vector (c_k, λ_k) . These variables help formulate the:

• inclusion constraint $\mathcal{Z}(c_k, G \cdot \operatorname{diag}(\lambda_k)) \subseteq U$:

$$s_i^{\top} c_k + \sum_j |s_i^{\top} G_j| \lambda_{jk} \le r_i, \quad \forall i,$$
 (11)

where $U = \{u : s_i^\top u \le r_i\}.$

• explicitly describe the volume⁴ $\operatorname{vol}(\mathcal{Z}(c_k, G\lambda_k))$:

$$\operatorname{vol}(\mathcal{Z}(c_k, G\Lambda_k)) = \sum_{1 \le j_1 < \dots j_n \le N} \left| \det(G^{j_1 \dots j_n}) \right| \cdot \prod_{j \in \{j_1 \dots j_n\}} \lambda_{jk}.$$
(12)

The formulation becomes simpler if the scaling is homogeneous $(\lambda_k^* = \lambda_{jk}, \forall j)$.



⁴Gover and Krikorian 2010.

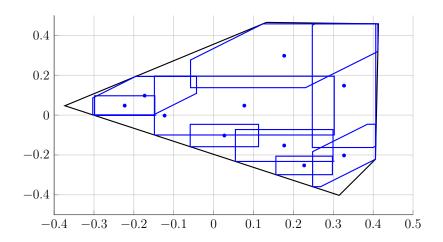
Implementation

We track the OMP formalism, without its theoretical convergence guarantees:

Algorithm 5: Area Coverage with zonotopic sets

- 1 Inputs: area to be covered $\it U$, sparsity constraint $\it s$
- 2 Result: pairs of centers and scaling factors (c_k, λ_k)
- 3 for k=1 to s do
- 4 Enlarge the zonotopes until they saturate the constraints
- Select Z_k where $k = \arg\min_k \operatorname{vol}(S_k \setminus Z_k)$
- 6 Update the uncovered area $\operatorname{vol}(S_{k+1}) = \operatorname{vol}(S_k \cup Z_k)$

Result



Thank You!

Questions?