Multi-Label Classification: Challenges, Pitfalls and Perspectives

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Joint work with Krzysztof Dembczynski, Willem Waegeman, Weiwei Cheng.



The distinguishing feature of multi-label classification (MLC): Instances may have **multiple** labels instead of only a **single** one!

X1	X2	X3	X4	Y1	Y2	Y3	Y4
0.34	0	10	174	0	0	1	0
1.45	0	32	277	0	1	0	0
1.22	1	46	421	0	0	0	1
0.74	1	25	165	0	1	0	0
0.95	1	72	273	1	0	0	0
1.04	0	33	158	0	0	1	0
0.92	1	81	382	?	?	?	?

In conventional classification, class labels are mutually exclusive.



The distinguishing feature of multi-label classification (MLC): Instances may have **multiple** labels instead of only a **single** one!

X1	X2	X3	X4	Y1	Y2	Y3	Y4
0.34	0	10	174	0	1	1	0
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0.92	1	81	382	?	?	?	?

Exploiting **label dependencies** (correlations) has become a major concern in MLC research.



Note that conventional multi-class classification is a special case of MLC, not the other way around ...



... and specific structure is normally exploited for restricted classes (e.g., linear programming).



The idea of having "additional information" comes from the "binary view" (which is closely related to multi-task and transfer learning): Multiple binary instead of a generalized multi-class problem.

X1	X2	X3	X4	English	French	Spanish	Dutch
0.34	0	10	174	0	1	1	0
1.45	0	32	277	0	1	0	1
1.22	1	46	421	0	0	0	1
0.74	1	25	165	0	1	1	1
0.95	1	72	273	1	0	1	0
1.04	0	33	158	1	1	1	0
0.92	1	81	382	?	?	?	?

Indeed, MLC problems can be "binarized" in an obvious way, and it would be solved if **binary relevance** (BR) learning was enough ...



The typical MLC paper:

- A new method is proposed that exploits label dependencies in one way or the other.
- It is shown to have better average accuracy than existing approaches in terms of a bunch of MLC loss functions.

Criticisms:

- Results are reported on average and not carefully analyzed: No investigation of the reasons for improvement and the conditions under which label dependencies are beneficial (some comments will follow).
- Implicitly, it seems to be assumed that one algorithm can be beneficial (if not optimal) for multiple loss functions (second part).
- The notion of "label correlation" is often used in a purely intuitive manner, without a precise formal definition (third part).













Adding noise by inverting each label with small probability (0.1).





The Label Powerset (LP) Approach



Treat every potential labeling as a new (meta-)class

ightarrow reduction to multi-class classification







Potential advantage of taking label dependence into account comes at the cost of more classes, imbalanced distributions, more complex decision boundaries.

Hamming and Subset 0/1 Loss





AGENDA



- 1. MLC Loss Functions and Risk Minimization
- 2. Label Dependence
- 3. Connections to Related Fields (maybe)
- 4. Concluding Remarks



- Given an instance space \mathcal{X} and label space $\mathcal{Y} = \{0, 1\}^m$, an MLC examples $(\boldsymbol{x}, \boldsymbol{y}) \in \mathcal{X} \times \mathcal{Y}$ is generated according to a probability distribution $\mathbf{P}(\mathbf{X}, \mathbf{Y})$.
- $\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y}) = \mathbf{P}(\mathbf{Y} | \boldsymbol{x})$ is the conditional distribution of \mathbf{Y} given $\mathbf{X} = \boldsymbol{x}$, and by $\mathbf{P}_{\boldsymbol{x}}^{(i)}(Y_i) = \mathbf{P}^{(i)}(Y_i | \boldsymbol{x})$ the corresponding marginal distribution of Y_i conditioned on \boldsymbol{x} :

$$\mathbf{P}_{x}^{(i)}(1) = \mathbf{P}^{(i)}(1 \,|\, x) = \sum_{y \in \mathcal{Y}: y_i = 1} \mathbf{P}(y \,|\, x)$$

- A multilabel classifier \mathbf{h} is an $\mathcal{X} \to \mathcal{Y}$ mapping that assigns a (predicted) labeling to each instance $x \in \mathcal{X}$. Thus, the output of a classifier \mathbf{h} is a vector

$$\mathbf{h}(\boldsymbol{x}) = (h_1(\boldsymbol{x}), h_2(\boldsymbol{x}), \dots, h_m(\boldsymbol{x})).$$

Example of a Conditional Probability



\mathbf{Y}_1	\mathbf{Y}_2	\mathbf{Y}_3	$\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y})$
0	0	0	0
0	0	1	0
0	1	0	0.4
1	0	0	0.3
0	1	1	0
1	0	1	0.3
1	1	0	0
1	1	1	0

Risk Minimization



The risk-minimizing model \mathbf{h}^* is defined in a pointwise way by

$$\mathbf{h}^*(\boldsymbol{x}) = \arg\min_{\boldsymbol{y}} \mathbb{E}_{\mathbf{Y}|\mathbf{X}} L(\mathbf{Y}, \boldsymbol{y}) ,$$

where $L(\cdot, \cdot)$ is a loss function defined on multi-label predictions.



The spectrum of losses in MLC is wider than in conventional classification!

• Subset 0/1 loss:

$$L_{0/1}(\boldsymbol{y}, \mathbf{h}(\boldsymbol{x})) = \llbracket \boldsymbol{y} \neq \mathbf{h}(\boldsymbol{x})
rbracket$$

• Hamming loss:

$$L_H(\boldsymbol{y}, \mathbf{h}(\boldsymbol{x})) = \frac{1}{m} \sum_{i=1}^m \llbracket y_i \neq h_i(\boldsymbol{x}) \rrbracket$$

• Rank loss:

$$L_{R}(\boldsymbol{y}, \mathbf{f}(\boldsymbol{x})) = \sum_{(i,j): y_{i} > y_{j}} \left(\left[\!\left[f_{i} < f_{j}\right]\!\right] + \frac{1}{2} \left[\!\left[f_{i} = f_{j}\right]\!\right] \right)$$

MLC Loss Functions



• For the subset 0/1 loss, the risk-minimizing prediction is given by the mode of the distribution:

$$\mathbf{h}^*(oldsymbol{x}) = rg\max_{oldsymbol{y}\in\mathcal{Y}} \mathbf{P}(oldsymbol{y}\,|\,oldsymbol{x})$$

• For the Hamming loss, the risk-minimizing prediction is given by

$$h_i^*(\boldsymbol{x}) = \arg \max_{b \in \{0,1\}} \mathbf{P}^{(i)}(b \,|\, \boldsymbol{x})$$

"joint mode" versus **"combined mode"** (combination of the modes of the margins)

Hamming Loss vs. Subset 0/1 Loss



\mathbf{Y}_1	\mathbf{Y}_2	\mathbf{Y}_3	$\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y})$
0	0	0	0
0	0	1	0
0	1	0	0.4
1	0	0	0.3
0	1	1	0
1	0	1	0.3
1	1	0	0
1	1	1	0

\mathbf{Y}_1	\mathbf{Y}_2	\mathbf{Y}_3	$\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y})$
0	0	0	0
0	0	1	0
0	1	0	0.4
1	0	0	0.3
0	1	1	0
1	0	1	0.3
1	1	0	0
1	1	1	0

subset 0/1 risk minimizer: (0, 1, 0)

would be the optimal choice of LP

Hamming risk minimizer: (1, 0, 0)

would be the optimal choice of BR



Define the regret of a classifier h with respect to a target loss L as follows:

$$r_L(\boldsymbol{h}) = R(\boldsymbol{h}) - R(\boldsymbol{h}^*)$$

= $\mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}}L(\boldsymbol{Y}, \boldsymbol{h}(\boldsymbol{X})) - \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}}L(\boldsymbol{Y}, \boldsymbol{h}^*(\boldsymbol{X})),$

with R the risk and h_z^* the Bayes-optimal classifier with respect to L.

Specifically for Hamming and subset 0/1 loss:

$$r_H(\boldsymbol{h}) = \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}} L_H(\boldsymbol{Y}, \boldsymbol{h}(\boldsymbol{X})) - \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}} L_H(\boldsymbol{Y}, \boldsymbol{h}_H^*(\boldsymbol{X}))$$
$$r_{0/1}(\boldsymbol{h}) = \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}} L_{0/1}(\boldsymbol{Y}, \boldsymbol{h}(\boldsymbol{X})) - \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}} L_{0/1}(\boldsymbol{Y}, \boldsymbol{h}_{0/1}^*(\boldsymbol{X}))$$



What is the "Hamming-regret" of a subset 0/1 optimal classifier and vice versa?

$$r_{H}(\boldsymbol{h}_{0/1}^{*}) = \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}}L_{H}(\boldsymbol{Y}, \boldsymbol{h}_{0/1}^{*}(\boldsymbol{X})) - \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}}L_{H}(\boldsymbol{Y}, \boldsymbol{h}_{H}^{*}(\boldsymbol{X}))$$

$$\uparrow$$

$$use classifier tailored for subset 0/1 Because (\mathbf{h}_{0/1}^{*}(\boldsymbol{X})) - \mathbb{E}_{\boldsymbol{X}\boldsymbol{Y}}L_{0/1}(\boldsymbol{Y}, \boldsymbol{h}_{0/1}^{*}(\boldsymbol{X}))$$

$$\uparrow$$

$$use classifier tailored for Hamming Subset 0/1 Because (\mathbf{h}_{0/1}^{*}(\boldsymbol{X}))$$



Proposition: The following upper bound holds:

$$\mathbb{E}_{\mathbf{Y}} L_{0/1}(\mathbf{Y}, \mathbf{h}_{H}^{*}(\mathbf{x})) - \mathbb{E}_{\mathbf{Y}} L_{0/1}(\mathbf{Y}, \mathbf{h}_{0/1}^{*}(\mathbf{x})) < 0.5.$$

Moreover, this bound is tight, i.e.,

$$\sup_{\mathbf{P}} \left(\mathbb{E}_{\mathbf{Y}} L_{0/1}(\mathbf{Y}, \boldsymbol{h}_{H}^{*}(\boldsymbol{x})) - \mathbb{E}_{\mathbf{Y}} L_{s}(\mathbf{Y}, \boldsymbol{h}_{0/1}^{*}(\boldsymbol{x})) \right) = 0.5,$$

where the supremum is taken over all probability distributions on \mathcal{Y} .

K. Dembczynski et al. *Regret Analysis for Performance Metrics in Multi-Label Classication: The Case of Hamming and Subset Zero-One Loss.* ECML-2010.



Proposition: The following upper bound holds for m > 3:

$$\mathbb{E}_{\boldsymbol{Y}}L_H(\boldsymbol{Y},\boldsymbol{h}_{0/1}^*(\boldsymbol{x})) - \mathbb{E}_{\boldsymbol{Y}}L_H(\boldsymbol{Y},\boldsymbol{h}_H^*(\boldsymbol{x})) < \frac{m-2}{m+2}.$$

Moreover, this bound is tight, i.e.

$$\sup_{\mathbf{P}}(\mathbb{E}_{\mathbf{Y}}L_H(\mathbf{Y}, \boldsymbol{h}_{0/1}^*(\boldsymbol{x})) - \mathbb{E}_{\mathbf{Y}}L_H(\mathbf{Y}, \boldsymbol{h}_H^*(\boldsymbol{x}))) = \frac{m-2}{m+2},$$

where the supremum is taken over all probability distributions on \mathcal{Y} .

A classifier tailored for subset 0/1 loss may perform extremely poor in terms of Hamming loss (at least theoretically)!

K. Dembczynski et al. *Regret Analysis for Performance Metrics in Multi-Label Classication: The Case of Hamming and Subset Zero-One Loss.* ECML-2010.



Empirically, we could confirm that BR is often better (!) than LP in terms of Hamming loss but worse in terms of subset 0/1 loss.

		BR SVM	BR MLRules	LPC pSVM	LPC+ pSVM	LPC SVM
	image	0.198 (4)	0.1928(2)	0.1888(1)	0.2021(5)	0.1954(3)
b0	scene	0.1071 (5)	0.0871(1)	0.0919(3)	0.095(4)	0.0891(2)
Ē	music	0.2049 (1)	0.208(2)	0.2091(3)	0.2232(5)	0.2119(4)
Ē	reuters	0.0663 (5)	0.0479(1)	0.0565(2)	0.0596(3)	0.0628(4)
2	yeast	0.2016 (1)	0.2086(3)	0.2156(4)	0.2523(5)	0.2075(2)
ц Т	genbase	0.0008 (1)	0.0015(5)	0.0011(3)	0.0012(4)	0.001(2)
	slashdog	0.048 (2)	0.0402(1)	0.0534(4)	0.0631(5)	0.0481(3)
	medical	0.0102 (1)	0.0106(2)	0.0132(4)	0.0135(5)	0.0115(3)
-	Average Rank	2.7	2.1	2.75	4.25	3.2

		BR SVM	BR MLRules	LPC pSVM	LPC SVM
	image	0.767 (4)	0.6705(3)	0.5595(2)	0.5315(1)
-	scene	0.4757 (4)	0.4221(3)	0.3299(2)	0.3008(1)
5	music	0.7538 (3)	0.7622(4)	0.7353(2)	0.6846(1)
	reuters	0.3735 (4)	0.2684(3)	0.2391(1)	0.2676(2)
S	yeast	0.8552 (3)	0.8643(4)	0.8155(2)	0.746(1)
D D	genbase	0.0211 (1)	0.0332(4)	0.0257(3)	0.0211(1)
S	slashdog	0.656 (2)	0.6721(3)	0.6819(4)	0.546(1)
	medical	0.3405 (2)	0.3497(3)	0.363(4)	0.3119(1)
	Average Rank	2.875	3.375	2.5	1.125

Regret Analysis

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Similar observations have been made for classification vs. ranking (e.g., learning AUC-optimizing classifiers):

A good classifier is not necessarily a good ranker!



Regret Analysis







- Even though the true differences will normally be smaller than suggested by these worst case estimates, or may even shrink to zero under specific assumptions (e.g., probability of mode > ½, conditional independence), different MLC loss functions will generally call for different classifiers.
- Stated differently, a single classifier is unlikely to be optimal for various loss functions at the same time.
- Thus, empirical studies suggesting the opposite should arouse suspicion ...



Train a label powerset classifier for each k-subset of labels, or a random subset thereof.

X1	X2	X3	X1		Y3	Y4
0.34	0	10	174		1	0
1.45	0	32	277		0	1
1.22	1	46	421		0	1
0.74	1	25	165		1	1
0.95	1	72	273		1	0
1.04	0	33	158		1	0
0.92	1	81	382		1	1

G. Tsoumakas and I. Vlahavas. *Random k-labelsets: An ensemble method for multilabel classication*. ECML 2007.

RAKEL: Prediction



(Y1, Y2) classifier \rightarrow	0
(Y1, Y3) classifier \rightarrow	1
(Y1, Y4) classifier \rightarrow	1

- (Y2, Y3) classifier \rightarrow
- (Y2, Y4) classifier \rightarrow
- (Y3, Y4) classifier \rightarrow

Y1	Y2	Y3	Y4
0	1		
1		1	
1			0
	0	1	
	0		1
		0	1
1	0	1	1

majority vote (or thresholding)

What is RAKEL Estimating?



\mathbf{Y}_1	\mathbf{Y}_2	\mathbf{Y}_3	\mathbf{Y}_4	$\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y})$
0	0	0	0	3/64
0	0	0	1	6/64
0	0	1	0	2/64
0	1	0	0	3/64
1	0	0	0	6/64
0	0	1	1	8/64
0	1	0	1	8/64
1	0	0	1	0
0	1	1	0	9/64
1	0	1	0	2/64
1	1	0	0	5/64
0	1	1	1	1/64
1	0	1	1	8/64
1	1	0	1	3/64
1	1	1	0	0
1	1	1	1	0

\mathbf{Y}_1	\mathbf{Y}_2	$\mathbf{P}_{oldsymbol{x}}(\mathbf{Y})$		
0 0 1 1	0 1 0 1	19/64 <mark>21/64</mark> 16/64 8/64	←	Bayes prediction for Rakel

Rakel prediction: (0, 0, 1, 1)

Hamming loss minimizer: (0, 0, 0, 1) subset 0/1 loss minimizer: (0, 1, 1, 0)



- For k=1, RAKEL equals **Binary Relevance** (hence tailored for Hamming).
- For k=m, RAKEL equals Label Powerset (hence tailored for subset 0/1).
- One may conjecture that, by varying k, RAKEL smoothly "interpolates" between Hamming and subset 0/1.
- Experiment for m=4, with probability distributions are selected uniformly at random.



AGENDA



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- 2. Label Dependence
- 3. Connections to Related Fields
- 4. Concluding Remarks

Risk Minimization



The risk-minimizing model \mathbf{h}^* is defined in a pointwise way by

$$\mathbf{h}^*(\boldsymbol{x}) = \arg\min_{\boldsymbol{y}} \mathbb{E}_{\mathbf{Y}|\mathbf{X}} L(\mathbf{Y}, \boldsymbol{y}) \; ,$$

where $L(\cdot, \cdot)$ is a loss function defined on multi-label predictions.

• Subset 0/1 loss:

$$\mathbf{h}^*(\boldsymbol{x}) = rg\max_{\boldsymbol{y}\in\mathcal{Y}}\mathbf{P}_{\boldsymbol{x}}(\boldsymbol{y})$$

• Hamming loss:

$$h_i^*(\boldsymbol{x}) = \arg \max_{b \in \{0,1\}} \mathbf{P}_{\boldsymbol{x}}^{(i)}(b)$$

Exploiting label dependence is probably more important for subset 0/1 than for Hamming!



Standard stochastic independence: The joint distribution is the product of the marginals.

Conditional distribution for an instance x form the (1,1) region, assuming independent error terms ϵ_1 and ϵ_2 :

$\mathbf{P}_{oldsymbol{x}}(\mathbf{Y})$	0	1	$\mathbf{P}_{\boldsymbol{x}}^{(1)}(Y_1)$
0	0.01	0.09	0.10
1	0.09	0.81	0.90
$\mathbf{P}_{\boldsymbol{x}}^{(1)}(Y_2)$	0.10	0.90	1

\rightarrow conditional label independence



Standard stochastic independence: The joint distribution is the product of the marginals.

Assuming a close dependency, namely $\epsilon_1 = \epsilon_2$:

$\mathbf{P}_{oldsymbol{x}}(\mathbf{Y})$	0	1	$\mathbf{P}_{\boldsymbol{x}}^{(1)}(Y_1)$
0	0.1	0	0.1
1	0	0.9	0.9
$\mathbf{P}_{\boldsymbol{x}}^{(1)}(Y_2)$	0.1	0.9	1

 \rightarrow conditional label dependence



- The conditional distribution is our target, as it allows for (Bayes) optimal prediction (regardless of the loss function).
- In other words, we can "exploit" this distribution, respectively the label dependencies it implies, for **optimal decision making**.
- However, by "exploiting label correlation" one normally means using them for learning, not for decision making.
- But how to exploit conditional label dependence from this point of view, given that it refers to a single instance?



Taking a global view on label correlation:

X2	X3	X1	Y1	Y2	Y3	Y4
0	10	174	0	1	1	0
0		277	0	1	0	1
1		421	0	0	0	1
			0	1	1	1
			1	0	1	0
			1	1	1	0
			0	1	1	1

Unconditional label dependence refers to the joint distribution $\mathbf{P}(\mathbf{Y})$ instead of $\mathbf{P}_{\boldsymbol{x}}(\mathbf{Y}) = \mathbf{P}(\mathbf{Y} \,|\, \boldsymbol{x}).$



$$\mathbf{P}(\mathbf{Y}) \,=\, \int_{\mathcal{X}} \mathbf{P}(\mathbf{Y} \,|\, oldsymbol{x}) \,d\mu(oldsymbol{x})$$

unconditional dependence = "average" conditional dependence

$$\begin{array}{c} \mathbf{P}(\mathbf{Y} \,|\, \mathbf{X}) \,\propto\, \mathbf{P}(\mathbf{X} \,|\, \mathbf{Y}) \times \mathbf{P}(\mathbf{Y}) \\ \uparrow & \uparrow \\ \\ \text{conditional} & \text{unconditional} \\ \end{array}$$

 \rightarrow exploiting unconditional dependence, typically via regularization, for predicting conditional distributions (or functions thereof)

Stacking



stacking classifier

$$(y_1, y_2) \mapsto (y_1^*, y_2^*)$$
 $(x_1, x_2) \mapsto (h_1(x_1, x_2), h_2(x_1, x_2)) = (y_1, y_2)$
base classifier (BR)

		pred	icted	tr	ue
		[]		[<u> </u>
x_1	x_2	y_1	y_2	y_1	y_2
0.34	-0.45	0	0	1	0
0.45	0.56	1	1	1	1
-0.22	0.82	0	0	0	1
0.74	-0.12	0	1	0	0

Stacking in the Running Example





Average loss as a function of the angle, with and without stacking



– base classifier = 1-NN

- stacking by table classifier

number of training examples = 50

Improvement on average!



Extended stacking classifier: $(x_1, x_2 \mid y_1, y_2) \mapsto (y_1^*, y_2^*)$





Extended stacking classifier: $(x_1, x_2 \mid y_1, y_2) \mapsto (y_1^*, y_2^*)$



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- Stacking, Bayesian inference and related regularization methods.
- Multi-task learning and transfer learning (also based on regularization)
- Reduced rank regression for multivariate regression with unconditionally dependent outputs.
- Label dimensionality reduction, e.g., kernel dependency estimation.
- Structured output prediction
- Conditional random fields and graphical models
- Preference Learning

Preference Learning





Preference Learning



Training	Prediction	
binary	binary	multi-label classification
binary	total order	multi-label classification (ranking)
graded	graded	graded muti-label classification (ICML 2010)
partial order	total order	label ranking
partial order	partial order	ranking with abstention (ECML 2010)

Calibrated Label Ranking



Combining absolute and relative evaluation:



F. Fürnkranz et al. Multilabel classification via calibrated label ranking. Machine Learning 73(2), 2008.



- Comparing MLC methods is difficult, since the effect of "taking label correlations into account" is hard to isolate from other changes that become necessary.
- MLC loss functions are of a quite different nature, and it's arguably impossible to minimize all of them by the same algorithm ...
- Label dependence can be considered at different levels (conditional, unconditional, in-between), and the concrete mechanisms of "taking label correlations into account" call for an explanation.
- MLC is related to many other subfields of ML (and statistics), and existing work in these fields should not be ignored.