# Wasserstein distance between Lévy measures with applications to Bayesian nonparametrics

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**Bocconi University** 



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## Joint work with:



Marta Catalano



Antonio Lijoi



Igor Prünster

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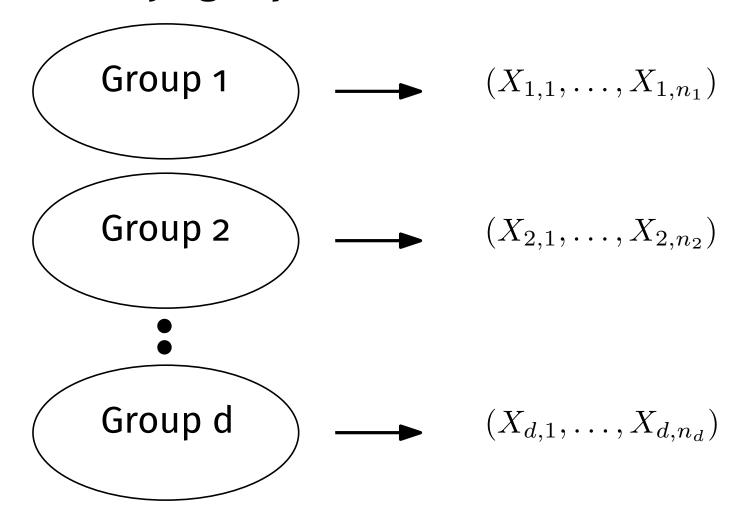
Igor Prünster

### **Disclaimer**

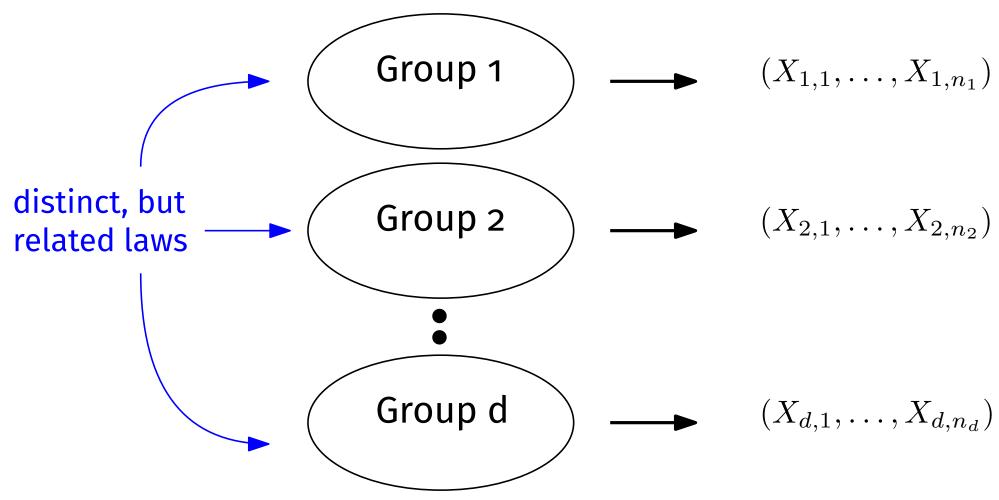
I am not a (Bayesian) statistican.

My background: mathematical analysis, optimal transport.

## **Quantifying dependence**

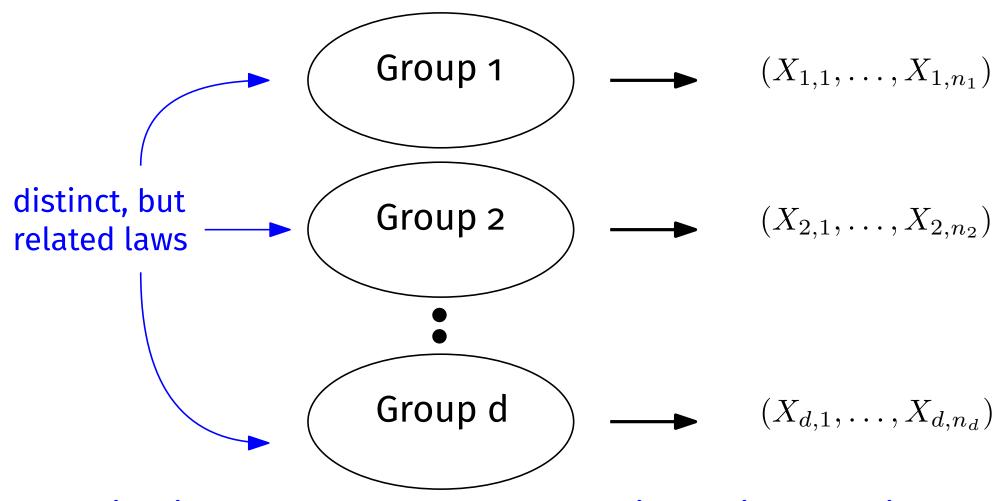


## **Quantifying dependence**



Bayesian inference allows for borrowing of information

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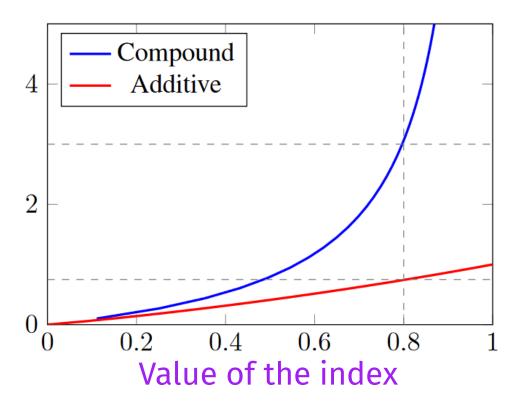
Bayesian inference allows for borrowing of information

**Goal**: quantifying the amount of **dependence** between groups already present in the **prior** 

### **Snapshot of the final result**

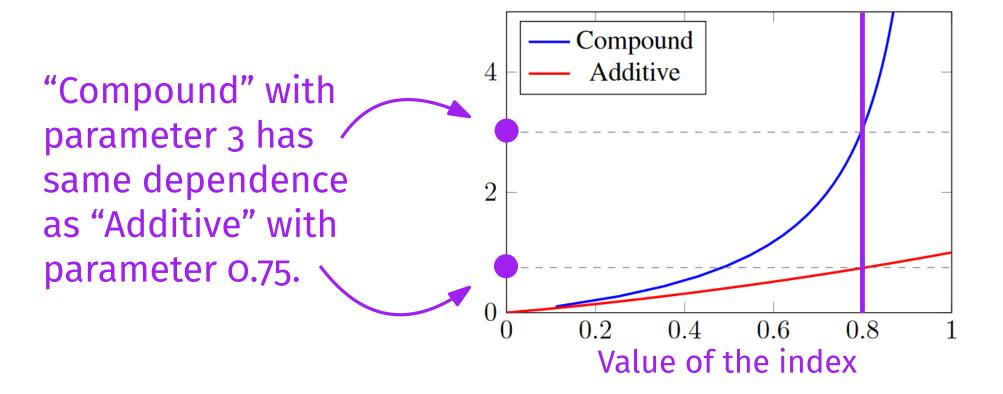
Our contribution: an index of dependence quantifying dependence in the prior

Different parametrized models of prior

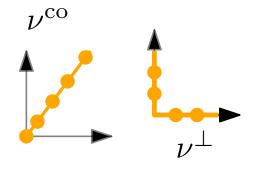


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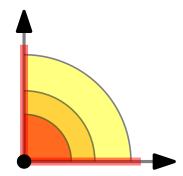


Allow for comparision between different priors



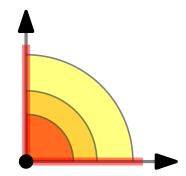
## 1 - Context, general strategy

## 2 - Building the index with optimal transport





## 2 - Building the index with optimal transport



## **Bayesian Non Parametrics**

 $ilde{p}$  random probability measure on  $\mathbb X$ 

$$X_1, X_2, \ldots, X_n | \tilde{p} \stackrel{\text{i.i.d.}}{\sim} \tilde{p}$$

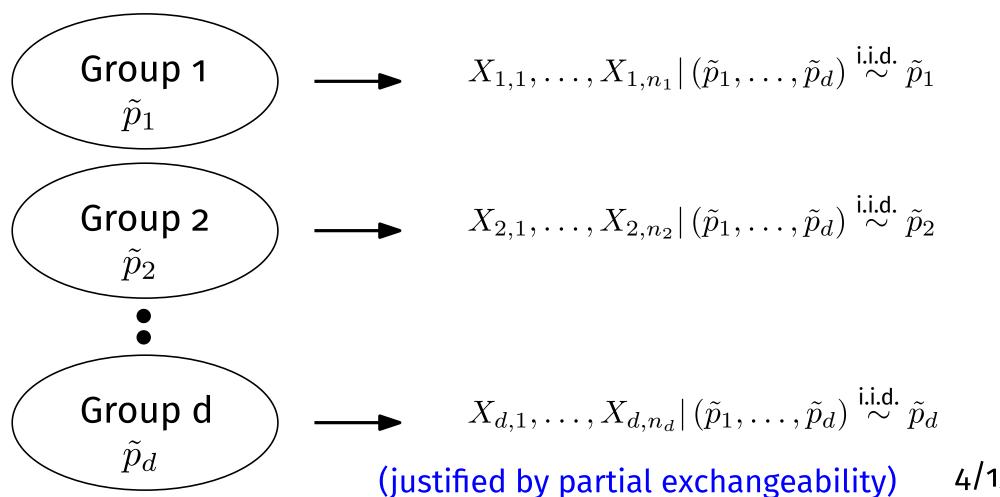
(justified by exchangeability)

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## **Specific setting: Completely Random Vectors**

$$\begin{split} \tilde{\boldsymbol{\mu}} &= (\tilde{\mu}_1, \tilde{\mu}_2, \dots, \tilde{\mu}_d) \text{ Completely Random Vector} \\ & X_{1,1}, X_{1,2}, \dots, X_{1,n_1} | \, \tilde{\boldsymbol{\mu}} \overset{\text{i.i.d.}}{\sim} \, \frac{\tilde{\mu}_1}{\tilde{\mu}_1(\mathbb{X})} \\ & X_{2,1}, X_{2,2}, \dots, X_{2,n_2} | \, \tilde{\boldsymbol{\mu}} \overset{\text{i.i.d.}}{\sim} \, \frac{\tilde{\mu}_2}{\tilde{\mu}_2(\mathbb{X})} \\ & \vdots \\ & X_{d,1}, X_{d,2}, \dots, X_{d,n_d} | \, \tilde{\boldsymbol{\mu}} \overset{\text{i.i.d.}}{\sim} \, \frac{\tilde{\mu}_d}{\tilde{\mu}_d(\mathbb{X})} \end{split}$$

**Definition (CRV).** For all  $A_1, \ldots, A_n \subseteq \mathbb{X}$  disjoints, the vectors  $\tilde{\boldsymbol{\mu}}(A_1), \ldots, \tilde{\boldsymbol{\mu}}(A_n)$  are independent random vectors in  $\mathbb{R}^d_+$ .

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Contains all dependence in the prior

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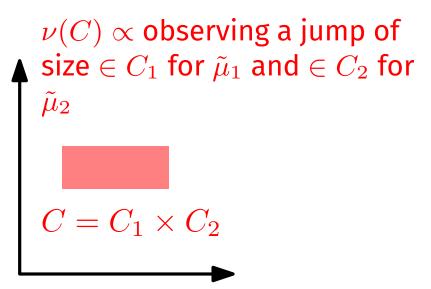
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For  $A \subseteq \mathbb{X}$ , the random variables  $\tilde{\mu}_1(A), \ldots, \tilde{\mu}_d(A)$  may be dependent.

Assumptions of **homogeneity** and no fixed atoms:

$$ilde{oldsymbol{\mu}} = \sum_{i=1}^{\infty} ilde{\mathbf{J}}_i \delta_{Y_i}$$

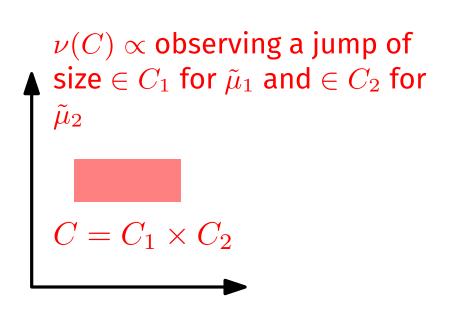
where  $(Y_i)_i \in \mathbb{X}$  (atoms) follow base measure  $P_0$ ; and  $(\tilde{\mathbf{J}}_i)_i$  (jumps) independent from  $(Y_i)_i$  follow Poisson point cloud on  $\mathbb{R}^d_+$  with intensity measure  $\nu$  (Lévy measure).

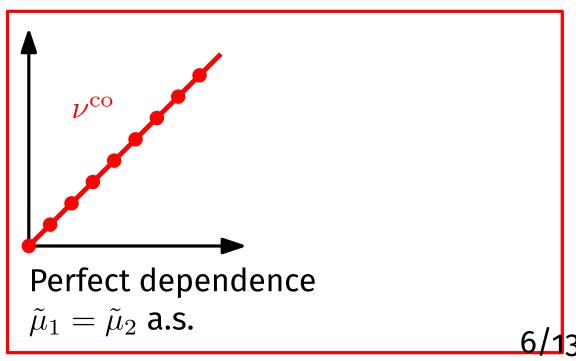


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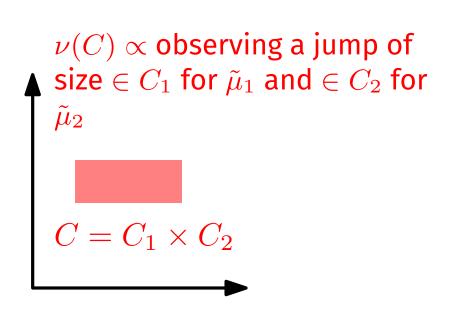


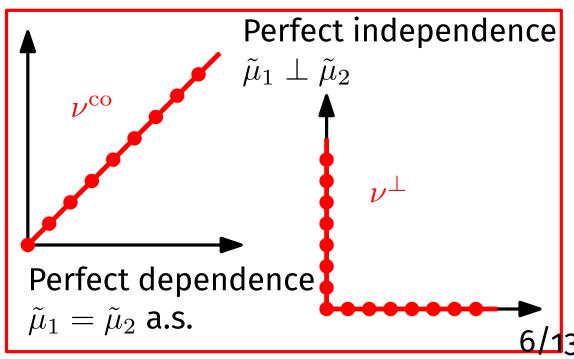


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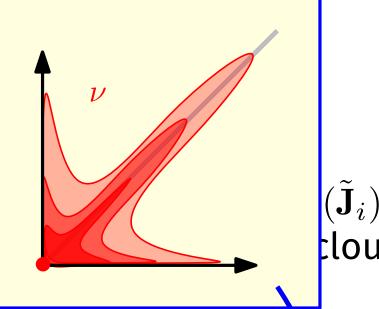




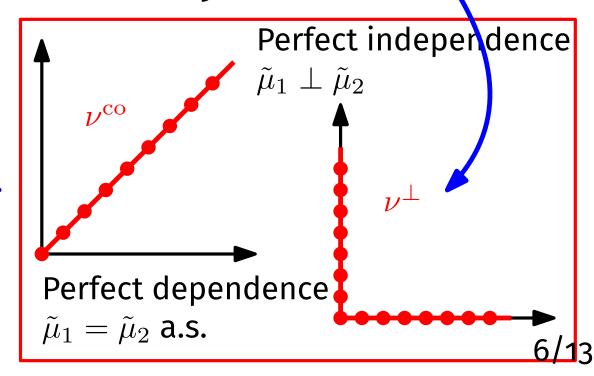
**Assumpt** 

where (Y (jumps) i on  $\mathbb{R}^d_+$  wi

**Goal**: distinguish between these two cases.



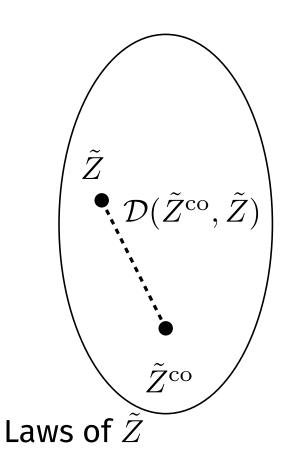
 $u(C) \propto ext{observing a jump of }$   $ext{size} \in C_1 ext{ for } ilde{\mu}_1 ext{ and } \in C_2 ext{ for }$   $ilde{\mu}_2$   $C = C_1 imes C_2$ 



## A general method to construct an index

### **Ingredients**:

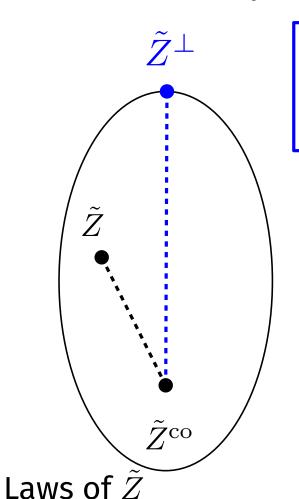
- $\tilde{Z}$  random object,  $\tilde{Z}^{\mathrm{co}}$  "most dependent".
- $\mathcal{D}$  "discrepancy" between laws of random objects.



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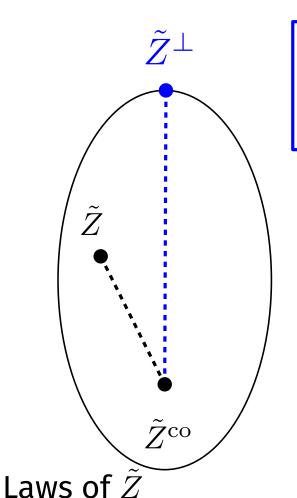


**To check**:  $\mathcal{D}(\tilde{Z}^{\mathrm{co}}, \tilde{Z})$  is maximized when  $\tilde{Z} = \tilde{Z}^{\perp}$  the most independent structure.

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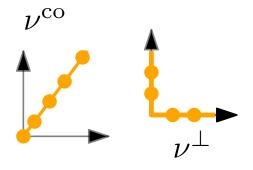
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#### Then define:

$$\mathcal{I}(\tilde{Z}) = 1 - \frac{\mathcal{D}(\tilde{Z}^{co}, \tilde{Z})}{\mathcal{D}(\tilde{Z}^{co}, \tilde{Z}^{\perp})}.$$

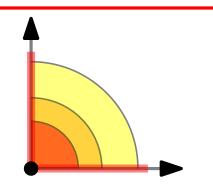
It belongs to [0, 1] and satisfies:

$$\mathcal{I}(\tilde{Z}^{\perp}) = 0, \qquad \mathcal{I}(\tilde{Z}^{co}) = 1.$$

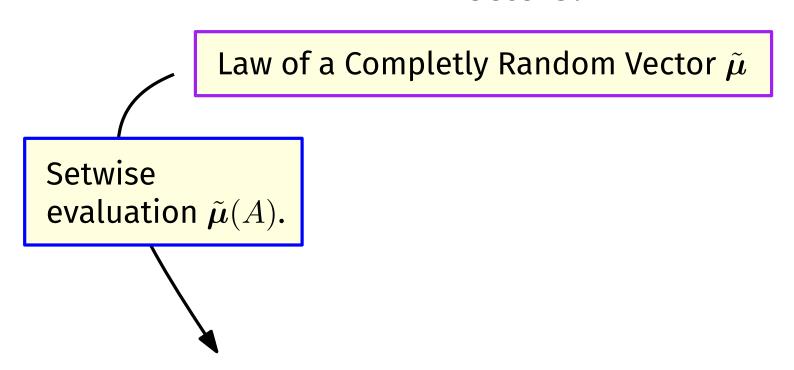


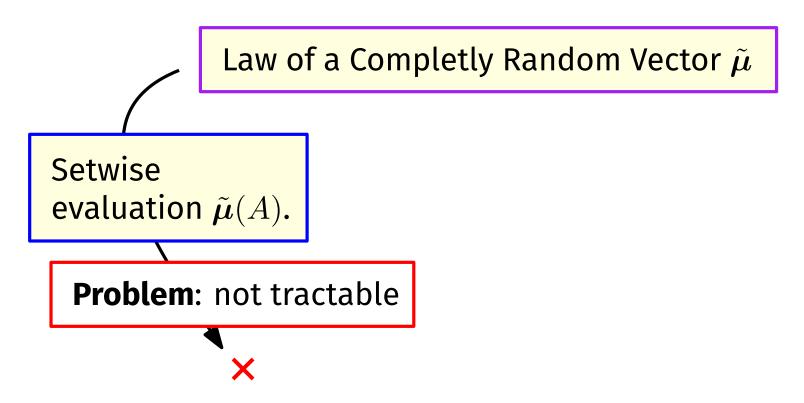
## 1 - Context, general strategy

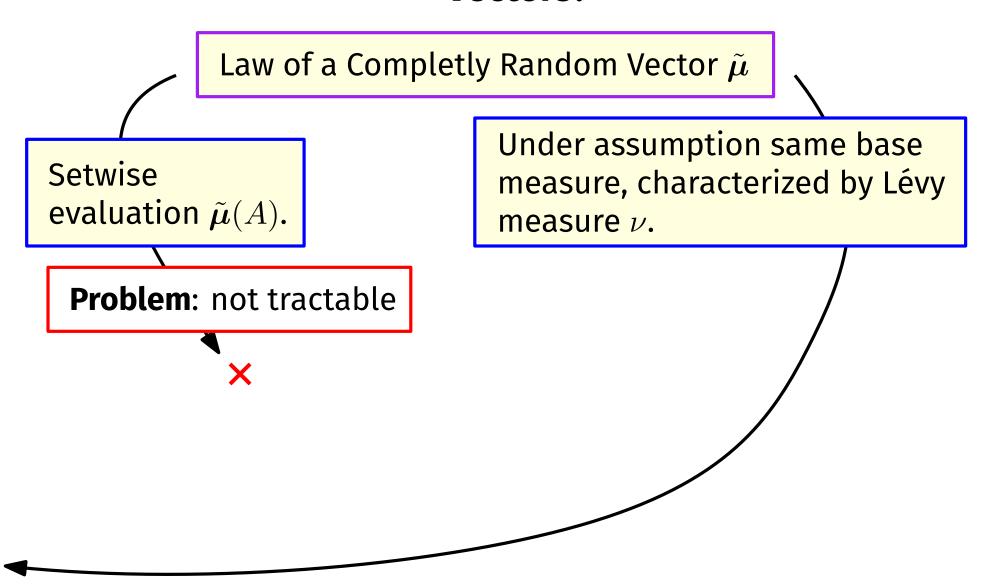
## 2 - Building the index with optimal transport

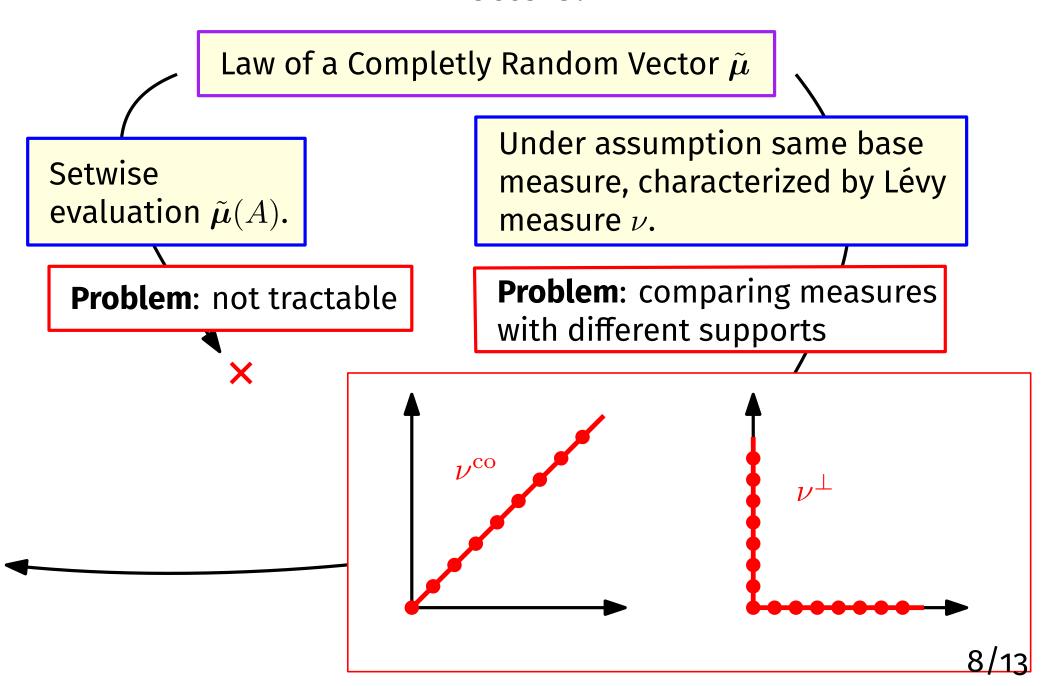


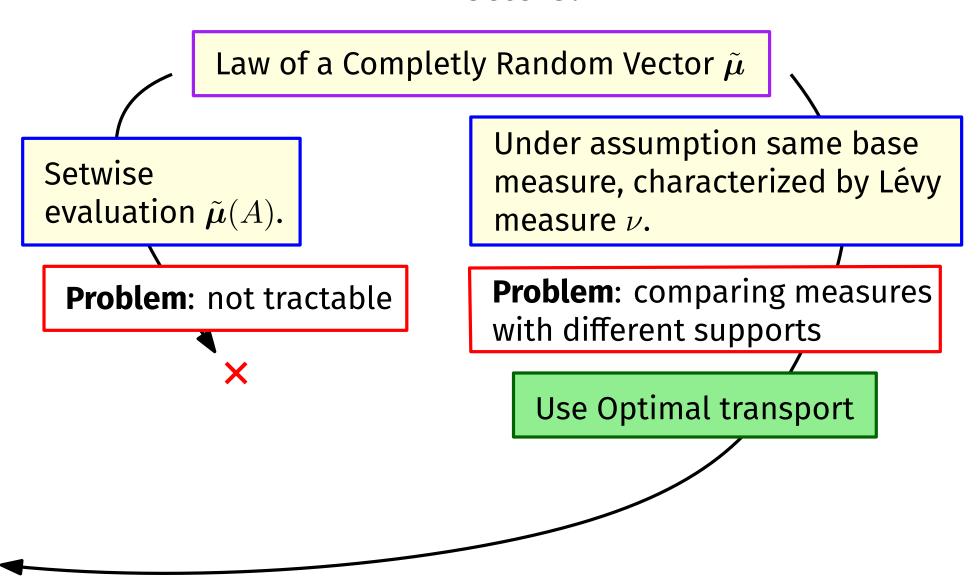
Law of a Completly Random Vector  $ilde{oldsymbol{\mu}}$ 

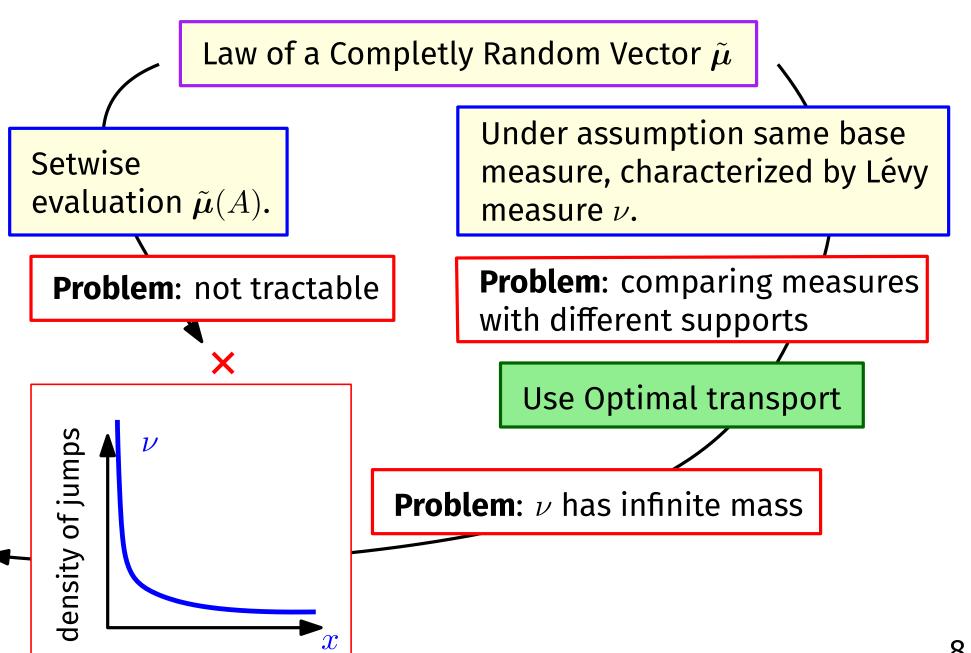


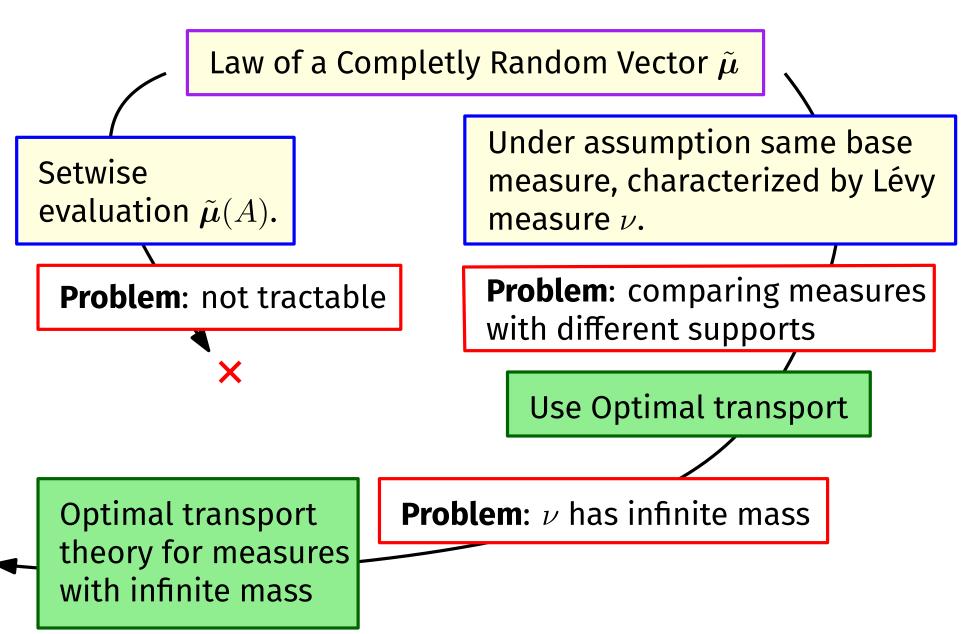


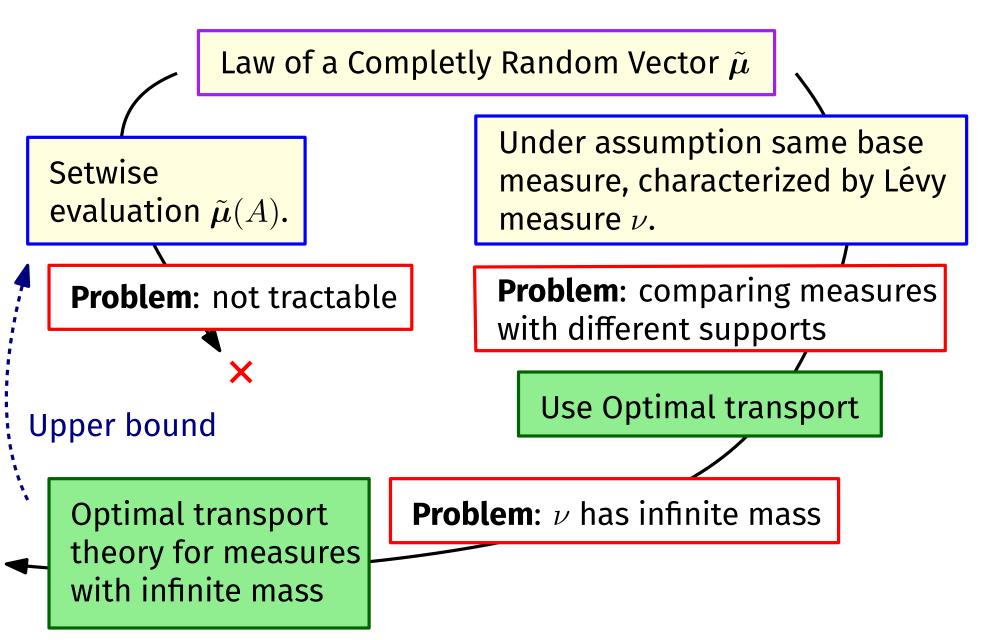












## (Classical) optimal transport

**Definition.** If  $\nu^1, \nu^2$  probability distributions, the Wasserstein distance is

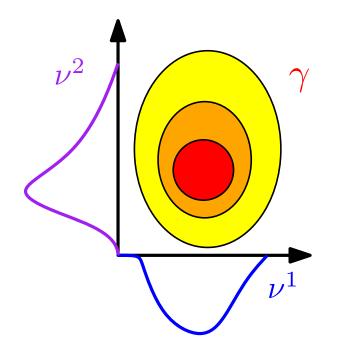
$$\mathcal{W}(\nu^1, \nu^2)^2 = \min_{(X,Y)} \left\{ \mathbb{E} \left[ \|X - Y\|^2 \right] \; : \; X \sim \nu^1 \; \text{and} \; Y \sim \nu^2 \right\}$$

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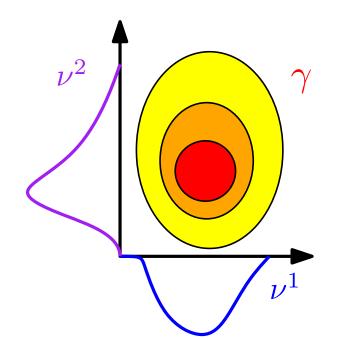


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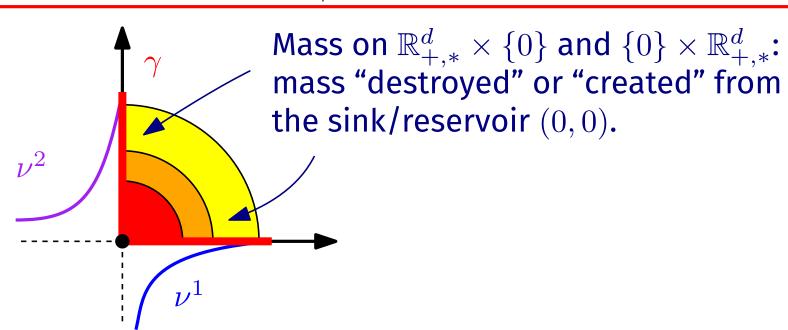
$$\leq \int ||x||^2 d\nu^1(x) + \int ||y||^2 d\nu^2(y)$$

**Observation**. Naively, makes sense if  $\nu^1, \nu^2$  have infinite mass but **finite** second moment.

#### **Extended Wasserstein distance**

**Definition**. If  $\nu^1, \nu^2$  positive measures on  $\mathbb{R}^d_+ \setminus \{0\}$  with **finite second moments**, the Wasserstein distance is

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 with  $\gamma$  measure on  $\mathbb{R}^{2d}_+ \setminus \{(0,0)\}$ .

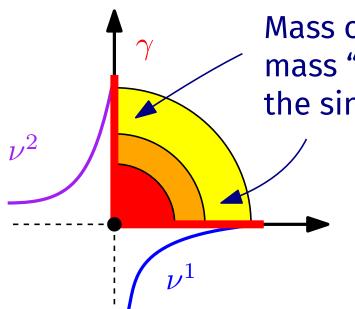


Figalli and Gigli (2010). A new transportation distance between non-negative measures, with applications to gradients flows with Dirichlet boundary conditions.

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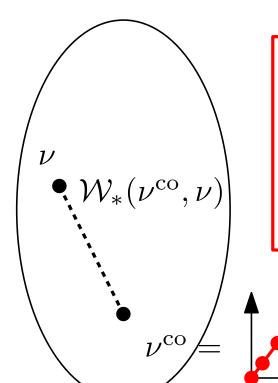
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 with  $\gamma$  measure on  $\mathbb{R}^{2d}_+ \backslash \{(0,0)\}$ .



Mass on  $\mathbb{R}^d_{+,*} \times \{0\}$  and  $\{0\} \times \mathbb{R}^d_{+,*}$ : mass "destroyed" or "created" from the sink/reservoir (0,0).

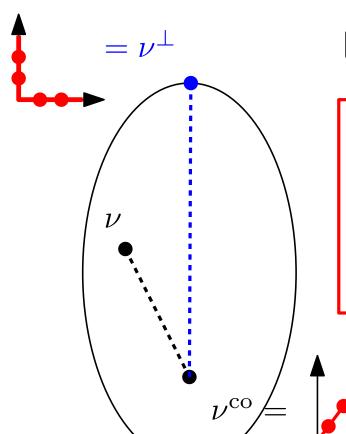
**In progress**: extending this idea to couple also the law of atoms for inhomogeneous CRV

Figalli and Gigli (2010). A new transportation distance between non-negative measures, with applications to gradients flows with Dirichlet boundary conditions.



First result.  $W_*(\nu^{co}, \nu)$  can be computed with 1d integrals of tail functions.

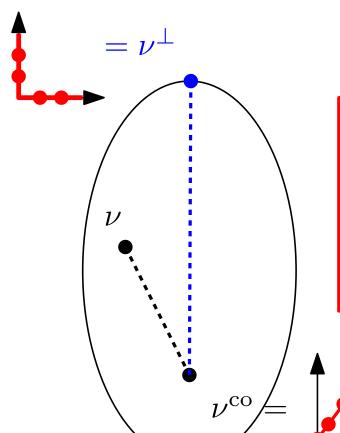
Space of Lévy measure over  $\mathbb{R}^d_+$  having same marginals



**First result**.  $\mathcal{W}_*(\nu^{\text{co}}, \nu)$  can be computed with 1d integrals of tail functions.

**Second result**. If  $\nu^{\text{co}}$  has infinite mass,  $\mathcal{W}_*(\nu^{\text{co}}, \nu)$  is maximized exactly for  $\nu = \nu^{\perp}$ .

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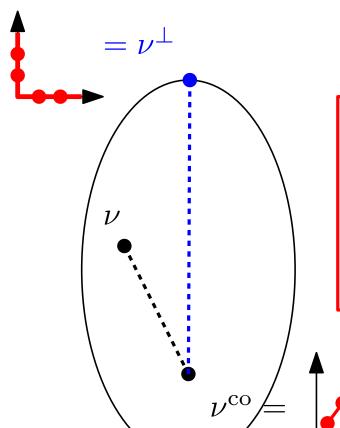
#### **Define:**

$$\mathcal{I}(\nu) = 1 - \frac{\mathcal{W}_*(\nu^{\text{co}}, \nu)^2}{\mathcal{W}_*(\nu^{\text{co}}, \nu^{\perp})^2}.$$

It belongs to [0,1] and satisfies:

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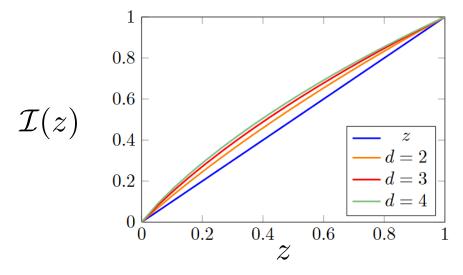
**Consequence**. We have an index of dependence for homogeneous infinitely active completely random vectors without fixed atoms, with equal marginals and finite second moments.

### **Examples**

### Additive model

Parameter  $z \in [0, 1]$ ,

$$\nu = (1 - z)\nu^{\perp} + z\nu^{\text{co}}$$



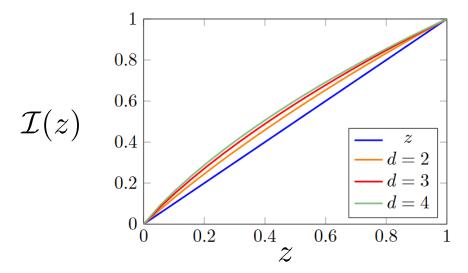
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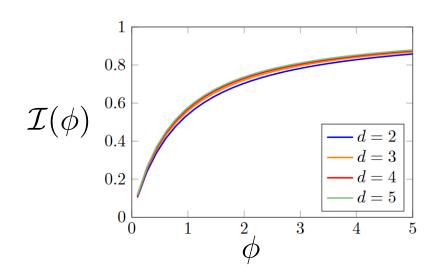
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### Compound random measures

Parameter  $\phi$  measures dependence

$$\nu(s_1, \dots, s_d) = \int_0^{+\infty} h^{\phi} \left(\frac{s_1}{u}, \dots, \frac{s_d}{u}\right) d\nu_*^{\phi}(u)$$

for well chosen  $h^{\phi}, \nu_*^{\phi}$ .



### **Examples**

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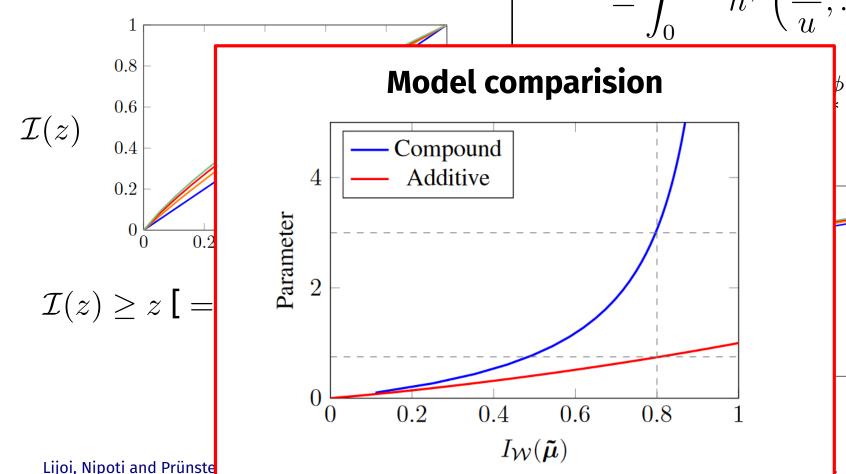
Griffin and Leisen (2017)

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12/13

d=2

d = 3d = 4

### **Conclusion**

#### What is done:

- Wasserstein distance between Lévy measures.
- Index of dependence between Completely Random Vectors.

#### What's next?:

- Study dependence in the posterior.
- Use this distance for other purposes: convergence of posterior, distance of a prior/posterior to a reference one, etc.

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Thank you for your attention