Estimation-of-Distribution Algorithms. Discrete Domain.

Petr Pošík

Dept. of CyberneticsČVUT FEL

[Introduction](#page-1-0) to EDAs

 Genetic [Algorithms](#page-2-0) andEpistasisGenetic [Algorithms](#page-3-0)

GA vs [EDA](#page-8-0)

[Content](#page-11-0) of the lectures

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Introduction to EDAs

Genetic Algorithms and Epistasis

From the lecture on epistasis: $x^{\text{best}} = 111...11, f(x^{\text{best}}) = 40$

GA works:

 $\boldsymbol{\nu}$ no dependencies

GA fails:

0

 0

0.10.20.3 0.4 0.5 0.6 0.70.80.91

xmean

f8x5bitTrap

- $\boldsymbol{\checkmark}$ deps. exist
- ✔ GA not able to work with them

Popsize160

0 5 10 15 20

bestaverage

generation

Popsize160

0 5 10_, 15 20

generation

GA works again:

- ✔deps. exist
- ✔GA knows about them

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A0M33EOA: Evolutionary Optimization Algorithms – ³ / ⁴⁴

Algorithm 1: Genetic Algorithm

7

- **Initialize** the population. **2**
- **3while** *termination criteria are not met* **do**
- 4 **Select** parents from the population. **4**
- **5EXECUTE: Cross over** the parents, create offspring.
- **6Mutate** offspring.
- **Incorporate** offspring into the population.

Select \rightarrow cross over \rightarrow mutate approach

- ✔are not adaptive, and
- ✔ cannot (or ususally do not) discover and use *the interactions among solution components*.

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What does an intearction mean?

- $\boldsymbol{\checkmark}$ we would like to create a new offspring by mutation
- ✔we would like the offspring to have better, or at least the same, quality as the paren^t
- $\boldsymbol{\mathsf{v}}$ if we must modify x_i together with x_j to reach the desired goal (if it is not possible to improve the solution by modifying either x_i or x_j only), then x_i interacts with x_j .

- ✔are not adaptive, and
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The goal of recombination operators:

 $\boldsymbol{\checkmark}$ Intensify the search in areas which contained "good" individuals in previous iterations.

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- $\boldsymbol{\checkmark}$ Intensify the search in areas which contained "good" individuals in previous iterations.
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The goal of recombination operators:

- ✔Intensify the search in areas which contained "good" individuals in previous iterations.
- ✔Must be able to take the interactions into account.
- ✔Why not directly describe the distribution of "good" individuals???
- $\boldsymbol{\nu}$ are not adaptive, and
- ✔ cannot (or ususally do not) discover and use *the interactions among solution components*.

GA vs EDA

Algorithm 1: Genetic Algorithm

Select \rightarrow cross over \rightarrow mutate approach

Algorithm 2: Estimation-of-Distribution Alg.

1 begin

45

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- **Initialize** the population. **2**
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- **Learn** a model of their distribution.
- **Sample** new individuals.

Incorporate offspring into the population.

Select \rightarrow model \rightarrow sample approach

GA vs EDA

Algorithm 1: Genetic Algorithm

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Explicit probabilistic model:

- ✔principled way of working with dependencies
- ✔adaptation ability (different behavior in different stages of evolution)

Algorithm 2: Estimation-of-Distribution Alg.

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Explicit probabilistic model:

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- ✔adaptation ability (different behavior in different stages of evolution)

Names:

- **EDA** Estimation-of-Distribution Algorithm
- **PMBGA** Probabilistic Model-Building Genetic Algorithm
- **IDEA** Iterated Density Estimation Algorithm

Select \rightarrow model \rightarrow sample approach

Content of the lectures

[Introduction](#page-1-0) to EDAs Genetic [Algorithms](#page-2-0) andEpistasis

Genetic [Algorithms](#page-3-0)

GA vs [EDA](#page-8-0)

[Content](#page-11-0) of the lectures

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

- 1. EDA for discrete domains (e.g. binary)
	- $\boldsymbol{\nu}$ Motivation example
	- ✔Without interactions
	- ✔Pairwise interactions
	- ✔Higher order interactions
- 2. EDA for real domain (vectors of real numbers)
	- ✔Evolution strategies
	- ✔Histograms
	- ✔Gaussian distribution and its mixtures

[Introduction](#page-1-0) to EDAs

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Motivation Example

Example

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)[Example](#page-13-0)

- Selection, [Modeling,](#page-14-0) Sampling
- UMDA [Behaviour](#page-15-0) for
- OneMax problem
- What about ^a [different](#page-16-0)fitness?
- UMDA behaviour on
- [concatanated](#page-17-0) traps
- What can be done about
- [traps?](#page-18-0)
- [Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

- EDAs without[interactions](#page-26-0)
- Pairwise [Interactions](#page-28-0)
- [Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

5-bit OneMax (CountOnes) problem:

- $\boldsymbol{\checkmark}$ f_{Dx1bitOneMax}(**x**) = $\sum_{d=1}^{D} x_d$
- \vee Optimum: 11111, fitness: 5

Algorithm: **Univariate Marginal Distribution Algorithm (UMDA)**

- $\boldsymbol{\nu}$ Population size: 6
- ✔ $\boldsymbol{\nu}$ Tournament selection: $t=2$
- $\boldsymbol{\nu}$ **Model:** vector of probabilities $p = (p_1, \ldots, p_D)$
	- ✘ each *^p^d* is the probability of observing ¹ at *^d*th element
- ✔ **Model learning:**
	- ✘ compute *^p* from selected individuals
- ✔ **Model sampling:**
	- ✘ generate ¹ on *^d*th position with probability *^p^d* (independently of other positions)

Selection, Modeling, Sampling

UMDA Behaviour for OneMax problem

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

- $\boldsymbol{\nu}$ 1s are better then 0s on average, selection increases the proportion of1s
- ✔ Recombination preserves and combines 1s, the ratio of 1s increases over time
- ✔ $\boldsymbol{\checkmark}$ If we have many 1s in population, we cannot miss the optimum

The number of evaluations needed for reliable convergence:

UMDA behaves similarly to GA with uniform crossover!

What about ^a different fitness?

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)[Example](#page-13-0) Selection, [Modeling,](#page-14-0) Sampling

 UMDA [Behaviour](#page-15-0) forOneMax problem

 What about ^a [different](#page-16-0)fitness?

 UMDA behaviour on[concatanated](#page-17-0) trapsWhat can be done about[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

For OneMax function:

✔UMDA works well, all the bits probably eventually converge to the right value.

Will UMDA be similarly successful for other fitness functions?

✔ Well, . no. :-(

Problem: **Concatanated 5-bit traps**

 $f = f_{\text{trap}}(x_1, x_2, x_3, x_4, x_5) +$ $+ f_{\text{trap}}(x_6, x_7, x_8, x_9, x_{10}) +$ $+$

The *trap* function is defined as

 $f_{\text{trap}}(\mathbf{x}) = \begin{cases} 5 & \text{if } u(\mathbf{x}) = 5 \\ 4 - u(\mathbf{x}) & \text{otherwise} \end{cases}$

where $u(\mathbf{x})$ is the so called *unity* function and returns the number of ¹^s in **^x** (it is actually the One Max function).

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)[Example](#page-13-0)

- Selection, [Modeling,](#page-14-0)
- Sampling
- UMDA [Behaviour](#page-15-0) for
- OneMax problem
- What about ^a [different](#page-16-0)fitness?

 UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about[traps?](#page-18-0)[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Traps:

- ✔Optimum in 111111...1
- ✔ \triangleright But *f*_{trap} (0 ∗ ∗ ∗ ∗) = 2 while *f*_{trap} (1 ∗ ∗ ∗ ∗) = 1.375
- ✔1-dimensional probabilities lead the GA to the wrong way!
- ✔ Exponentially increasing population size is needed, otherwise GA will not find optimum reliably.

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Example](#page-13-0)

Selection, [Modeling,](#page-14-0)

Sampling

UMDA [Behaviour](#page-15-0) for

OneMax problem

What about ^a [different](#page-16-0)

fitness?

 UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

The *f*trap function is *deceptive*:

- ✔Statistics over 1**** and 0**** do not lead us to the right solution
- ✔The same holds for statistics over $11***$ and $00***$, $111**$ and $000**$, $1111*$ and 0000*

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Example](#page-13-0)

Selection, [Modeling,](#page-14-0)

Sampling

UMDA [Behaviour](#page-15-0) for

OneMax problem

 What about ^a [different](#page-16-0)fitness?

 UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about

[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

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- ✔ Harder than the *needle-in-the-haystack* problem:
	- \boldsymbol{x} regular haystack simply does not provide any information, where to search for the needle
	- ✘ *f*trap-haystack actively lies to you—it points you to the wrong par^t of the haystack

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Example](#page-13-0)

- Selection, [Modeling,](#page-14-0)
- Sampling
- UMDA [Behaviour](#page-15-0) for
- OneMax problem
- What about ^a [different](#page-16-0)fitness?
- UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

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[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Example](#page-13-0)

- Selection, [Modeling,](#page-14-0)
- Sampling
- UMDA [Behaviour](#page-15-0) for
- OneMax problem
- What about ^a [different](#page-16-0)fitness?
- UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

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Model learning:

- ✔build model for each 5-tuple of bits
- $\boldsymbol{\checkmark}$ compute $p(00000)$, $p(00001)$, ..., $p(11111)$,

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Example](#page-13-0)

- Selection, [Modeling,](#page-14-0)
- Sampling
- UMDA [Behaviour](#page-15-0) for
- OneMax problem
- What about ^a [different](#page-16-0)fitness?
- UMDA behaviour on[concatanated](#page-17-0) traps

What can be done about[traps?](#page-18-0)

[Good](#page-23-0) news!

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

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Model learning:

- $\boldsymbol{\checkmark}$ build model for each 5-tuple of bits
- $\boldsymbol{\checkmark}$ compute $p(00000)$, $p(00001)$, ..., $p(11111)$,

Model sampling:

- ✔Each 5-tuple of bits is generated independently
- ✔Generate 00000 with probability $p(00000)$, 00001 with probability $p(00001)$, ...

Good news!

Good statistics work great!

What shall we do next?

If we were able to

- ✔ find goo^d statistics with ^a small overhead, and
- ✔use them in the UMDA framework,

we would be able to solve order-*k* separableproblems using $\mathcal{O}(D^2)$ evaluations.

 $\boldsymbol{\checkmark}$... and there are many problems of this type.

The problem solution is closely related to the so-called *linkage learning*, i.e. discovering and using statistical dependencies among variables. [Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

 Discrete EDAs: **[Overview](#page-25-0)**

 EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Discrete EDAs

Discrete EDAs: Overview

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs Discrete EDAs: **[Overview](#page-25-0)**

 EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

- 1. Overview:
	- (a) Univariate models (without interactions)
	- (b) Bivariate models (pairwise dependencies)
	- (c) Multivariate models (higher order interactions)
- 2. Conclusions

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

 EDAs without[interactions](#page-27-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

EDAs without interactions

EDAs without interactions

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0) EDAs without[interactions](#page-27-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

- 1. **Population-based incremental learning (PBIL)**Baluja, 1994
- 2. **Univariate marginal distribution algorithm (UMDA)**Mühlenbein and Paaß, 1996
- 3. **Compact genetic algorithm (cGA)**Harik, Lobo, Goldberg, 1998

Similarities:

 $\boldsymbol{\checkmark}$ all of them use a vector of probabilities

Differences:

- ✔ PBIL and cGA do not use population (only the vector *^p*); UMDA does
- \triangleright PBIL and cGA use different rules for the adaptation of *p*

Advantages:

- $\boldsymbol{\mathsf{v}}$ Simplicity
- ✔ Speed
- ✔ Simple simulation of large populations

Limitations:

✔ Solves reliably only order-1 decomposable problems

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise

[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

EDAs with Pairwise Interactions

From single bits to pairwise models

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0) dependency tree

 Example of [dependency](#page-33-0)tree learning

[Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

How to describe two positions together?

 $\boldsymbol{\checkmark}$ Using the joint probability distribution:

Number of free parameters:

✔Using statistical dependence:

Number of free parameters:

 $p(A, B) = p(B|A) \cdot p(A)$: *p*(*^B* =1|*A*= $p(B = 1 | A = 0)$
 $p(B = 1 | A = 1)$ *p*(*^A*=1)

Question: what is the number of parameters in case of the following models?

AB $\mathsf C$ AB $\mathsf C$ AB $\mathsf C$

From single bits to pairwise models

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0) dependency tree

 Example of [dependency](#page-33-0)tree learning

[Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

How to describe two positions together?

 $\boldsymbol{\checkmark}$ Using the joint probability distribution:

Number of free parameters: 3

✔Using statistical dependence:

Number of free parameters:

 $p(A, B) = p(B|A) \cdot p(A)$: *p*(*^B* =1|*A*= $p(B = 1 | A = 0)$
 $p(B = 1 | A = 1)$ *p*(*^A*=1)

Question: what is the number of parameters in case of the following models?

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[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0) dependency tree

 Example of [dependency](#page-33-0)tree learning

[Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

How to describe two positions together?

 $\boldsymbol{\checkmark}$ Using the joint probability distribution:

Number of free parameters: 3

✔Using statistical dependence:

Number of free parameters: 3

 $p(A, B) = p(B|A) \cdot p(A)$: *p*(*^B* =1|*A*= $p(B = 1 | A = 0)$
 $p(B = 1 | A = 1)$ *p*(*^A*=1)

Question: what is the number of parameters in case of the following models?

AB $\mathsf C$ AB $\mathsf C$ AB $\mathsf C$

Example with pairwise dependencies: dependency tree

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0) dependency tree

Example of [dependency](#page-33-0)tree learning

[Dependency](#page-39-0) tree: probabilities

EDAs with pairwise

[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

- $\boldsymbol{\checkmark}$ Nodes: binary variables (loci of chromozome)
- ✔Edges: dependencies among variables
- ✔ Features:
	- ✘Each node depends at most on ¹ other node
	- ✘Graph does not contain cycles
	- ✘Graph is connected

Learning the structure of dependency tree:

1. Score the edges using mutual information:

$$
I(X,Y) = \sum_{x,y} p(x,y) \cdot \log \frac{p(x,y)}{p(x)p(y)}
$$

- 2. Use any algorithm to determine the maximum spanning tree of the graph, e.g. Prim (1957)
	- (a) Start building the tree from any node
	- (b) Add such ^a node that is connected to the tree by the edge with maximumscore

 $\overline{8}$

 X_3

 x_2

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency treeExample of [dependency](#page-33-0)

tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency treeExample of [dependency](#page-33-0)

tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

EDAs with pairwise interactions

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

- 1. **MIMIC** (sequences)
	- $\boldsymbol{\nu}$ Mutual Information Maximization for Input**Clustering**
	- ✔de Bonet et al., ¹⁹⁹⁶
- 2. **COMIT** (trees)
	- ✔ Combining Optimizers with Mutual Information Trees
	- ✔Baluja and Davies, ¹⁹⁹⁷
- 3. **BMDA** (forrest)
	- $\boldsymbol{\nu}$ Bivariate Marginal Distribution Algorithm
	- \vee Pelikan and Mühlenbein, 1998

Summary

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)From [single](#page-29-0) bits topairwise models

Example with pairwise[dependencies:](#page-32-0)

dependency tree

 Example of [dependency](#page-33-0)tree learning

 [Dependency](#page-39-0) tree: probabilities

EDAs with pairwise[interactions](#page-45-0)

[Summary](#page-46-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

- ✔ Advantages:
	- ✘Still simple
	- ✘Still fast
	- ✘Can learn *something* about the structure
- ✔ Limitations:
	- ✘Reliably solves only order-2 decomposable problems

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[ECGA](#page-48-0) ECGA: [Evaluation](#page-50-0)metric [BOA](#page-51-0) BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

EDAs with Multivariate Interactions

ECGA

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[ECGA](#page-48-0)

 ECGA: [Evaluation](#page-50-0)metric[BOA](#page-51-0) BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Extended Compact GA, Harik, 1999

Marginal Product Model (MPM)

- ✔Variables are treated in groups
- ✔Variables in different groups are considered statistically independent
- ✔Each group is modeled by its joint probability distribution
- ✔The algorithm adaptively searches for the groups during evolution

ECGA

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[ECGA](#page-48-0)

 ECGA: [Evaluation](#page-50-0)metric[BOA](#page-51-0) BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

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- ✔The algorithm adaptively searches for the groups during evolution

Learning the structure

- 1. Evaluation metric: Minimum Description Length (MDL)
- 2. Search procedure: greedy
	- (a) Start with each variable belonging to its own group
	- (b) Perform such ^a join of two groups which improves the score best
	- (c) Finish if no join improves the score

ECGA: Evaluation metric

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[ECGA](#page-48-0)

 ECGA: [Evaluation](#page-50-0)metric

[BOA](#page-51-0) BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Minimum description length:

Minimize the number of bits needed to store the model and the data encoded using themodel

DL(*Model*, *Data*) ⁼ *DLModel* ⁺ *DLData*

Model description length:

Each group *g* has $|g|$ dimensions, i.e. $2^{|g|} - 1$ frequencies, each of them can take on values up to *^N*

 $DL_{Model} = \log N \sum_{g \in G}$ $(2^{|g|} - 1)$

Data description length using the model:

Defined using the entropy of marginal distributions (X_g is $|g|$ -dimensional random vector, x_g is its realization):

$$
DL_{Data} = N \sum_{g \in G} h(X_g) = -N \sum_{g \in G} \sum_{x_g} p(X_g = x_g) \log p(X_g = x_g)
$$

BOA

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions**[ECGA](#page-48-0)** ECGA: [Evaluation](#page-50-0)metric

[BOA](#page-51-0)

 BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

Bayesian Optimization Algorithm: Pelikán, Goldberg, Cantù-Paz, 1999 Bayesian network (BN)

- ✔Conditional dependencies (instead groups)
- ✔Sequence, tree, forrest — special cases of BN
- ✔For trap function:

- ✔ The same model used independently in
	- $\boldsymbol{\mathsf{x}}$ Estimation of Bayesian Network Alg. (EBNA), Etxeberria et al., ¹⁹⁹⁹
	- \boldsymbol{X} Learning Factorized Density Alg. (LFDA), Mühlenbein et al., 1999

BOA: Learning the structure

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[ECGA](#page-48-0) ECGA: [Evaluation](#page-50-0)metric

[BOA](#page-51-0)

 BOA: [Learning](#page-52-0) thestructure

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

- 1. Evaluation metric:
	- ✔Bayesian-Dirichlet metric, or
	- ✔Bayesian information criterion (BIC)
- 2. Search procedure: greedy
	- (a) Start with graph with no edges (univariate marginal product model)
	- (b) Perform one of the following operations, choose the one which improves thescore best
		- $\boldsymbol{\nu}$ Add an edge
		- $\boldsymbol{\nu}$ Delete an edge
		- ✔Reverse an edge
	- (c) Finish if no operation improves the score

BOA solves order- k decomposable problems in less then $\mathcal{O}(D^2)$ evaluations!

 $n_{evals} = \mathcal{O}(D^{1.55})$ to $\mathcal{O}(D^2)$

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)Test [function](#page-55-0) (cont.)[Scalability](#page-56-0) analysis[OneMax](#page-58-0) [Non-dec.](#page-59-0) Equal Pairs [Decomp.](#page-60-0) Equal Pairs Non-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) TrapModel structure during

[Conclusions](#page-66-0)

[evolution](#page-64-0)

Scalability Analysis

Test functions

One Max:

$$
f_{Dx1bitOneMax}(\mathbf{x}) = \sum_{d=1}^{D} x_d
$$

Equal Pairs:

$f_{\text{DbitEqualPairs}}(\mathbf{x}) = 1 +$ *D* ∑*d*=2 $f_{\text{EqualPair}}(x_{d-1}, x_d)$ *f*_{EqualPair} (x_1, x_2) =

Trap:

$$
f_{D\text{bitTrap}}(\mathbf{x}) = \begin{cases} D & \text{if } u(\mathbf{x}) = D \\ D - 1 - u(\mathbf{x}) & \text{otherwise} \end{cases}
$$

$$
f_{\text{EqualPair}}(x_1, x_2) = \begin{cases} 1 & \text{if } x_1 = x_2 \\ 0 & \text{if } x_1 \neq x_2 \end{cases}
$$

Sliding XOR:

$$
f_{\text{DbitSlidingXOR}}(\mathbf{x}) = 1 + f_{\text{AllEqual}}(\mathbf{x}) + f_{\text{AllEqual}}(\mathbf{x}) + f_{\text{AllEqual}}(\mathbf{x}) = \begin{cases} 1 & \text{if } \mathbf{x} = (000 \dots 0) \\ 1 & \text{if } \mathbf{x} = (111 \dots 1) \\ 0 & \text{otherwise} \end{cases} + \sum_{d=3}^{D} f_{\text{XOR}}(x_{d-2}, x_{d-1}, x_d)
$$
\n
$$
f_{\text{XOR}}(x_1, x_2, x_3) = \begin{cases} 1 & \text{if } x_1 \oplus x_2 = x_3 \\ 0 & \text{otherwise} \end{cases}
$$

Concatenated short basis functions:

$$
f_{NxKbitBasisFunction} = \sum_{k=1}^{K} f_{BasisFunction}(x_{K(k-1)+1}, \ldots, x_{Kk})
$$

P. Pošík © 2010

Test function (cont.)

- 1. *f*40x1bitOneMax
	- $\boldsymbol{\nu}$ order-1 decomposable function, no interactions
- 2. *f*1x40bitEqualPairs
	- $\boldsymbol{\nu}$ non-decomposable function
	- $\boldsymbol{\checkmark}$ weak interactions: optimal setting of each bit depends on the value of the preceding bit
- 3. *f*8x5bitEqualPairs
	- $\boldsymbol{\nu}$ order-5 decomposable function
- 4. *f*1x40bitSlidingXOR
	- $\mathbf v$ non-decomposable function
	- $\boldsymbol{\checkmark}$ stronger interactions: optimal setting of each bit depends on the value of the 2 preceding bits
- 5. *f*8x5bitSlidingXOR
	- \vee order-5 decomposable function
- 6. *f*8x5bitTrap
	- $\boldsymbol{\nu}$ order-5 decomposable function
	- $\boldsymbol{\checkmark}$ interactions in each 5-bit block are very strong, the basis function is deceptive

Scalability analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs[Decomp.](#page-60-0) Equal PairsNon-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

[Conclusions](#page-66-0)

Facts:

- $\boldsymbol{\checkmark}$ using small population size, population-based optimizers can solve only easy problems
- ✔ increasing the population size, the optimizers can solve increasingly harder problems
- ✔... but using ^a too big population is wasting of resources.

Scalability analysis

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs[Decomp.](#page-60-0) Equal PairsNon-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

[Conclusions](#page-66-0)

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- $\boldsymbol{\checkmark}$... but using a too big population is wasting of resources.

Scalability analysis:

- $\boldsymbol{\nu}$ determines the optimal (smallest) population size, with which the algorithm solves the given problem reliably
	- ✘reliably: algorithm finds the optimum in ²⁴ out of ²⁵ runs)
	- ✘ for each problem complexity, the optimal population size is determined e.g. using the bisection method
- ✔ studies the influence of the problem complexity (dimensionality) on the optimal population size and on the number of needed evaluations

Scalability on the One Max function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs [Decomp.](#page-60-0) Equal Pairs Non-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

Scalability on the non-decomposable Equal Pairs function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs [Decomp.](#page-60-0) Equal Pairs Non-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

Scalability on the decomposable Equal Pairs function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

[Non-dec.](#page-59-0) Equal Pairs

[Decomp.](#page-60-0) Equal Pairs

 Non-dec. [Sliding](#page-61-0) XOR [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap

 Model structure during[evolution](#page-64-0)

Scalability on the non-decomposable Sliding XOR function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

[Non-dec.](#page-59-0) Equal Pairs

[Decomp.](#page-60-0) Equal Pairs

Non-dec. [Sliding](#page-61-0) XOR

 [Decomp.](#page-62-0) Sliding XOR[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

Scalability on the decomposable Sliding XOR function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs[Decomp.](#page-60-0) Equal Pairs

Non-dec. [Sliding](#page-61-0) XOR

[Decomp.](#page-62-0) Sliding XOR

[Decomp.](#page-63-0) Trap Model structure during[evolution](#page-64-0)

Scalability on the decomposable Trap function

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

 [Non-dec.](#page-59-0) Equal Pairs[Decomp.](#page-60-0) Equal Pairs

Non-dec. [Sliding](#page-61-0) XOR

[Decomp.](#page-62-0) Sliding XOR

[Decomp.](#page-63-0) Trap

 Model structure during[evolution](#page-64-0)

Model structure during evolution

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

Test [functions](#page-54-0)

Test [function](#page-55-0) (cont.)

[Scalability](#page-56-0) analysis

[OneMax](#page-58-0)

[Non-dec.](#page-59-0) Equal Pairs

[Decomp.](#page-60-0) Equal Pairs

Non-dec. [Sliding](#page-61-0) XOR

[Decomp.](#page-62-0) Sliding XOR

[Decomp.](#page-63-0) Trap

 Model structure during[evolution](#page-64-0)

[Conclusions](#page-66-0)

During the evolution, the model structure is increasingly precise and at the end of theevolution, the model structure describes the problem structure exactly.

Model structure during evolution

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

- [Scalability](#page-53-0) AnalysisTest [functions](#page-54-0)Test [function](#page-55-0) (cont.)
- [Scalability](#page-56-0) analysis
- [OneMax](#page-58-0)
- [Non-dec.](#page-59-0) Equal Pairs[Decomp.](#page-60-0) Equal Pairs
- Non-dec. [Sliding](#page-61-0) XOR
- [Decomp.](#page-62-0) Sliding XOR

[Decomp.](#page-63-0) Trap

 Model structure during[evolution](#page-64-0)

[Conclusions](#page-66-0)

During the evolution, the model structure is increasingly precise and at the end of theevolution, the model structure describes the problem structure exactly.

NO! That's not true!

Why?

- \triangleright In the beginning, the distribution patterns are not very discernible, models similar to uniform distributions are used.
- \triangleright In the end, the population converges and contains many copies of the same individual (or ^a few individuals). No interactions among variables can be learned. Model structure is wrong (all bits independent), but the model describes theposition of optimum very precisely.
- \checkmark The model with the best matching structure is found somewhere in the middle of the evolution.
- ✔ Even though the right structure is never found during the evolution, the problemcan be solved successfully.

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

[Summary](#page-67-0)

 [Suggestions](#page-68-0) for discreteEDAs

Summary

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

[Summary](#page-67-0)

 [Suggestions](#page-68-0) for discreteEDAs

Models:

- ✔Bayesian networks are genera^l models of joint probability
- ✔High-dimensional models are hard to train
- $\boldsymbol{\checkmark}$ High-dimensional models are very flexible

Advantages:

✔Reliably solves problems decomposable to subproblems of bounded order

Limitations:

✔ Does not solve problems decomposable to logarithmic subproblems (hierarchical problems)

Suggestions for discrete EDAs

[Introduction](#page-1-0) to EDAs

Motivation [Example](#page-12-0)

[Discrete](#page-24-0) EDAs

EDAs without[interactions](#page-26-0)

Pairwise [Interactions](#page-28-0)

[Multivariate](#page-47-0) Interactions

[Scalability](#page-53-0) Analysis

[Conclusions](#page-66-0)

[Summary](#page-67-0)

 [Suggestions](#page-68-0) for discreteEDAs

For simple problems:

- $\boldsymbol{\checkmark}$ PBIL, UMDA, cGA
- $\boldsymbol{\checkmark}$ they behave similarly to simple GAs

For harder problems:

- $\boldsymbol{\checkmark}$ MIMIC, COMIT, BMDA
- $\boldsymbol{\nu}$ they are able to account for bivariate dependencies

For hard problems:

- ✔BOA, ECGA, EBNA, LFDA
- $\boldsymbol{\checkmark}$ they can take into account more general dependencies, problems with hierarchichal structures

For even harder problems:

 $\boldsymbol{\nu}$ hBOA (hierarchical BOA)