
Generative Modeling With Vine Models

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1 Formulation

The goal is to optimize the construction of the vine copula model in order to minimize the Wasserstein distance between the true underlying distribution and the distribution of model. The construction of the first tree can be seen as a sequential decision making process: in each step we select a variable and link it with a variable in the existing tree. Therefore it is natural to model the process in the syntax of Reinforcement Learning.

Let \mathbb{P}_r be the true underlying distribution and \mathbb{P}_g be the distribution of the vine copula model used to generated synthetic data. The Wasserstein-1 distance is defined as:

$$W(\mathbb{P}_r, \mathbb{P}_g) = \sup_{||f||_L \leq 1} \mathbb{E}_{x \sim \mathbb{P}_r}[f(x)] - \mathbb{E}_{x \sim \mathbb{P}_g}[f(x)]$$

where the supremum is over all the 1-Lipschitz functions. There are two choices that can be made:

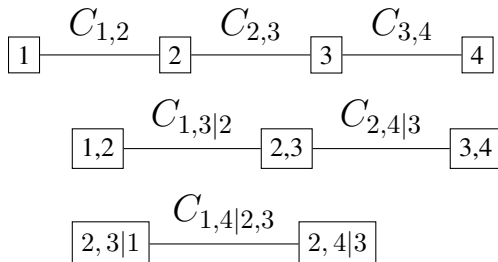
1. The choices of bi-variate copula functions (Is that possible to locally search for the best bi-variate copula function subject to some constraints.)
2. The sequence of variables to join when construct the vine

The construction of the first tree can be seen as a sequential decision making process: in each step we select a variable and link it to a variable in the existing tree.

For a fix generator, the discriminator solves the following optimizatin problem:

2 Related Work

3 Sampling



23 The sampling procedure requires a complete vine model of n nodes X_1, X_2, \dots, X_n , and their
 24 corresponding marginal distribution F_1, F_2, \dots, F_n . Considering the case where we have a D-Vine
 25 model with 4 nodes. We start by sampling uni-variate distribution u_1, u_2, u_3, u_4 from Uniform
 26 Distribution over $[0,1]$. We randomly pick a node to start with, say X_2 .

27 Then the first variable $x_2 \sim X_2$ can be sampled as:

$$x_2 = F_2^{-1}(u_2) \quad (1)$$

28 After we have x_2 , we randomly pick a node connected to X_2 . Suppose we pick X_3 , recall that the
 29 conditional density $f_{3|2}$ can be written as:

$$f_{3|2}(x_3|x_2) = f_3(x_3)c_{2,3}(F_2(x_2), F_3(x_3)) \quad (2)$$

$$= f_3(x_3)c_{2,3}(u_2, u_3) \quad (3)$$

$$= f_3(x_3)c_{3|2}(u_3) \quad (4)$$

30 Thus, $x_3 \sim X_3|X_2$ can be sampled by:

$$x_3 = F_3^{-1}(C_{3|2}^{-1}(u_3)) \quad (5)$$

31 where $C_{3|2}$ can be obtained from $C_{2,3}$ in T_1 by plugging in sampled values of u_2 .

32 Similarly, we pick a node that shares an edge with X_3 , say X_4 . Then $x_4 \sim X_4|X_2, X_3$ can be
 33 sampled as:

$$x_4 = F_4^{-1} \circ C_{4|3}^{-1} \circ C_{4|23}^{-1}(u_4) \quad (6)$$

34 Finally $x_1 \sim X_1|X_2, X_3, X_4$ can be sampled as:

$$x_1 = F_1^{-1} \circ C_{1|4}^{-1} \circ C_{1|34}^{-1} \circ C_{1|234}^{-1}(u_1) \quad (7)$$

35 For a more general vine graph, we can use a modified Breadth First Search to traverse the tree
 and keep track of nodes that have already been sampled. The general procedure is described below:

Algorithm 1 Generative Modeling with Vine Models

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1: procedure SAMPLING
2:   explore  $\leftarrow$  Queue()
3:   visited  $\leftarrow$  list()
4:   start  $\leftarrow$  randomly chosen start node
5:   explore.enqueue(start)
6:   while explore not empty do
7:     cur  $\leftarrow$  explore.dequeue()
8:     i  $\leftarrow$  len(visited)
9:      $x_{cur} = F_{cur}^{-1} \circ C_{cur|visited[i-1]}^{-1} \cdots \circ C_{cur|visited[1, \dots, i-1]}^{-1}(u_{cur})$ 
10:    for s  $\in$  neighbor(cur) do
11:      if s  $\in$  visited then
12:        Continue
13:      else
14:        explore.enqueue(s)
15:        visited.left_append(cur)

```

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37 4 Experiment

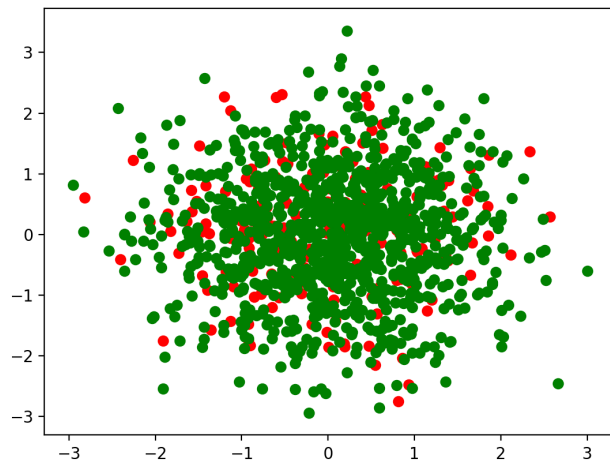


Figure 1: Independent Gaussian

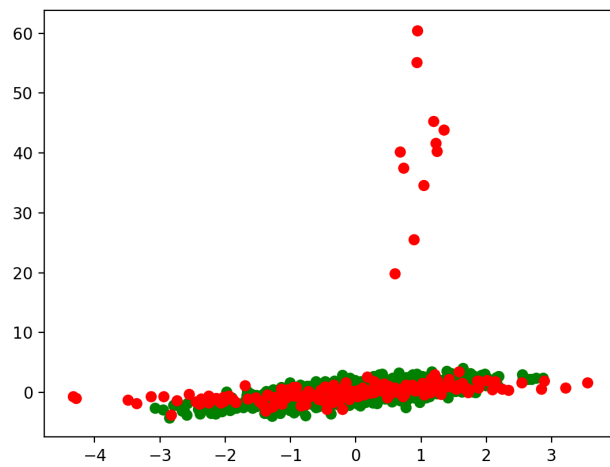


Figure 2: Correlated Gaussian