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Featured Application: This work will be applied to editable image generation. In the future, controllable 3D models can be generated to reduce the workload of film and television or game workers.

Abstract: Generative adversarial networks (GANs) are often used to generate realistic images, and GANs are effective in fitting high-dimensional probability distributions. However, during training, they often produce model collapse, which is the inability of the generative model to map the input noise to the real data distribution. In this work, we propose a model for disentanglement and mitigating model collapse inspired by the relationship between Hessian and Jacobian matrices. This is a concise framework for producing few modifications to the original model while facilitating the disentanglement. Compared to the pre-improvement generative models, our approach modifies the original model architecture only marginally and does not change the training method. Our method shows consistent resistance to model collapse on some image datasets, while outperforming the pre-improvement method in terms of disentanglement.

Keywords: deep learning; unsupervised learning; disentanglement; generative adversarial networks

1. Introduction

In recent years, many frameworks for representing high-dimensional data in an unsupervised manner have proliferated. Among many deep generative methods, generative adversarial networks (GANs) are among the most prominent methods for synthesizing real images. Generative Adversarial Network (GAN) has many meaningful works in high-resolution graph generation [1–5]. The emergence of GANs has made a huge leap in modeling high-dimensional distributed data, which has attracted great research interest. It consists of two networks: a generator and a discriminator, where the generator tries to map the latent data to the real data distribution. However, these data are often highly diverse. Therefore, the training of datasets often brings various model training collapse, both in terms of reduced model generation quality and generation diversity.

To solve this problem, there are three main options: (1) modifying the model architecture to change different architectures for different situations (e.g., AdaGAN [6], D2GAN [7]); (2) achieving better data mapping for the required generative goals by modifying the loss function (e.g., WGAN [8], LSGAN [9], EBAGN [10]); or (3) modifying the hidden space (DeLiGAN [11]). We focus on the model architecture, and we modify the original learning paradigm. This approach does not require large modifications to the original model, which greatly broadens the ease of use and applicability, and allows faster convergence to an equilibrium point.

Such a modified approach makes the model more entangled in the latent space. To cope with this problem, SeFa [12] disassembled the first fully connected layer of the GAN to disentanglement, and it is this layer of disentanglement that makes the overall disentanglement capability very limited. The Hessian GAN [13] minimizes the non-diagonal



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Copyright: © 2022 by the authors. Licensee MDPI, Basel, Switzerland. This article is an open access article distributed under the terms and conditions of the Creative Commons Attribution (CC BY) license (https:// creativecommons.org/licenses/by/ 4.0/). terms of the Hessian matrix to better learn disentanglement, and this method treats each output independently but lacks holistic constraint, so it lacks competitiveness in spatial representation. To solve this problem, Jacobian (OroJaR) [14] computes the Jacobian matrix of outputs to represent the changes caused by potential inputs, constraining the Jacobian vectors of each dimension to be orthogonal, so it can make it possible to compensate for this problem. Still, this method lacks the hessian penalty for effective access to the second-order of the original model information and weakens the output diversity. We demonstrate the intrinsic connection between these two methods and propose a new method that allows the model to have a clear separation boundary and effective spatial separation.

The mode collapse of the generative adversarial network has existed since the model was produced. Some scholars' works [15–18] make the model better to avoid this problem. For using these methods [15–18], the original model needs to be significantly changed, so we propose our method to mitigate mode collapse.

The method we propose to mitigate mode collapse is dedicated to further reducing the probability of mode collapse based on the original model, which we mainly show through loss line graphs and evaluation metrics. Moreover, our model can mitigate mode collapse without increasing the training burden.

Alleviating model collapse is only the first step in practical training. To generate higher-quality and controllable images, disentanglement is necessary. Some scholars' works [19–23] have done many works, but they do not perform well in disentanglement with multiple objects and complex scenes, so we propose our disentanglement method. Our method outperforms previous disentanglement performances for multiple objects and complex scenes.

In summary, we propose a new training framework that improves model stability and disentanglement with brief modifications to the model. Our approach allows training the model without getting into a local Nash equilibrium problem and doing the model as a disentanglement process. Our approach encourages discriminators and generators to diversify the generated samples to step out of the local optimum. At the same time, the loss term can make the boundaries of the diverse samples clearer to achieve the diversity of real data. In this framework that we proposed, we can choose to change the model architecture to resist model collapse, or we can choose to add our loss function to solve the entanglement. We conducted some experiments on various GAN models to evaluate the performance of our approach under different datasets. Our method consistently outperforms the preimprovement learning paradigm in model collapse mitigation and generation quality. In short, this paper makes the following contributions:

- We analyze the intrinsic connections and differences between Hessian regularization and Jacobian regularization.
- We propose a joint Hessian–Jacobian regularization model that enables GAN to encourage models to better disentanglement, while modifying the network architecture using only the hyperparameters in it can also better mitigate model collapse.
- We have done relevant experiments to show that our approach is useful in both mitigating model collapse and disentanglement representations.

2. Related Work

In deep learning, the two main types of unsupervised disentanglement learning are adversarial generative network (GAN) [24] and variational autoencoder (VAE) [25]. Before this, reconstruction networks already existed [26,27], and the creation of GANs brought at the same time the related optimization and disentanglement problems. Among the responses, there are changes in different architectures to modify the model architecture (AdaGAN [6], D2GAN [7]), and by modifying the loss function to achieve better data mapping for the required generative target (WGAN [8], LSGAN [9], EBAGN [10]), and by modifying the hidden space (DeLiGAN [11]).

2.1. Better Model Structure

2.1.1. Modifying the Model Architecture

In the training of the datasets, not all their data streams are connected. So, there are cases where a single generator has difficulty capturing all the streams. Some researchers have turned their attention to the network architecture of GAN. The main focus here is to modify the number of generators or discriminators. AdaGAN [6] improves model performance by aggregating many potential individual predictors to form a composite predictor, but its multiple models and complex training process make training more expensive. BEGAN [28] uses a self-encoder as a classifier and matches the loss distribution of the self-encoder through the loss based on Wasserstein distance. In order to make 3D controllable, Henzler et al. [29] used differentiable rendering learning to represent voxels, and the voxels based on this method were not clear enough to cause artifacts. Heusel et al. [30] proposed two time-scale update rules. Still, this approach tends to make the training of GANs network hard. Even if the regularization term is added, the effect is not very obvious, and the training cost is also high. CGAN [31] is different from GAN, it adds the same condition to both the Generator and the Discriminator and uses this condition to determine the direction of the Generation model. CGAN [31] uses the conditional probability Generation model to generate multiple outputs simultaneously through different inputs of dependent variables. D2GAN [7] uses two discriminators to minimize KL divergence (Kullback-Leibler divergence: KL divergence is used to express the degree of difference between two distributions) and invert KL divergence, and it can solve the model collapse problem to some extent. However, like other methods, D2GAN [7] has the common problem of the computational complexity.

2.1.2. Alternative Loss Functions

InfoGAN [32] proposed an information-theoretic extension of GANs that obtains a deconvoluted representation of the data by penalized latent reconstructions. Mao et al. [9] proposed LSGAN, which uses a least-squares loss function instead of the GAN loss function, and they defined a pullback operator to map the generated samples to the data stream shape, which alleviates the instability of GAN training and poor quality of the generated images. EBGAN [10] provides an energy-based explanation for GAN, and thus can use a series of energy-based tools for GAN. The exhaustive grid search experiment verifies the complete set of hyperparameters and architecture settings of GAN and EBGAN [10]. Radford et al. [33] alleviates the stability of GAN-generated images to some extent, but the generated images lack global uniformity. Arjovsky et al. [8] used Wasserstein distance instead of JS scatter while completing stable training, which is difficult to train and slow to converge during practical experiments. Then, Gulrajani et al. [34] improved WGAN [8] and proposed WGAN-GP, which used weight clipping in dealing with Lipschitz constraints. The method 3D Controllable Image Synthesis proposed by Liao et al. [35] requires pure background images as an auxiliary to process multi-object scenes, which brings differentiable rendering to a new level, and the difficulty of obtaining pure background images reduces to a certain extent the breadth of application of this method. Hessian Penalty [13] achieved orthogonalization by constraining the output to be diagonal to the input Hessian matrix. After that, OroJaR [14] constrained the input by Jacobian matrix dimensions to disentangle the model. BGAN [36] means Boundary Seeking GAN, and it is dedicated to finding the boundary of GAN; this refers to the boundary when the discriminator's loss is generally stable at 0.5 because the effect of generating pictures at this time is the best. To obtain the optimal *D*, Hjelm et al. [36] made some modifications to the loss function. When D(x) = 0.5, G is optimal. The opposite approach has been used by [37], in order to make localization operations easier and more meaningful in each semantic region, they used a large number of localization controls by semantic mappings and sorted them, thus achieving the desired effect.

2.2. Disentanglement

Among the existing unsupervised methodological approaches, the two main types of disentanglement representations for GAN are disentanglement in the GAN potential space and regularized disentanglement. GANSPACE [38] applies PCA to the sampled data to find significant and meaningful directions in the style space of StyleGAN [1]. Mildenhall et al. [39] proposed a neural radiation field that combines implicit neural models with volume rendering to reconstruct complex scenes. FineGAN [40] uses information theory with weaker supervised information for background, object pose, shape, and texture to find meaningful orientations. An unsupervised method such as [41] enables the classifier to correctly identify the semantic orientations in the matrix. InfoGAN [32] includes autoencoders on the potential space; however, it has been shown to have similar stability issues to standard GAN and requires stabilization of empirical skills. Giraffe [42] advocates generation directly in the 3D modeling process, and they introduce regularization in 3D modeling. Giraffe [42] can better disentangle a single object, but the disentanglement does not perform well in the multi-object case.

In generation tests, generators still suffer from uneven generation distributions, including collapsed generation content and incomplete semantics. However, these works provide a variety of approaches and solutions that alleviate the generator crash problem to some extent, but the disentanglement does not perform well. Among many existing generative frameworks, GANs tend to synthesize the highest quality generation. However, they are more difficult to optimize due to unstable training dynamics. Compared to previous methods, our approach requires neither too much change in the model structure nor any additional trainable parameters to achieve a better Nash equilibrium and less overfitting, and a better disentanglement representation.

3. Theoretical Method and Simplified Calculation

In this section, we first introduce the basic theory of GAN dynamics description. Then, we describe the intrinsic unity of the proposed Hessian regularization and Jacobian regularization for learning the disentanglement representation. Finally, we discuss it with related methods (Jacobian (OroJaR) [14] and Hessian penalty [13]) and present our approach.

3.1. Basic Theory and Dynamics of GAN

Generative adversarial networks (GANs) fit the data distribution of generators to the original dataset using a game in which generative models compete with discriminators. The generator G takes a random noise vector as input and maps it to points in the target data distribution. The discriminator D receives the data and tries to determine whether it really comes from the empirical distribution (if so, it outputs 1) or is made by the generator (output 0). During the training iterations, noise vectors from the Gaussian distribution G are pushed through the generator network to form a batch of generated data samples. This process is expressed in the max-min formula as Equation (1)

$$\min_{u} \max_{v} \mathcal{L}(u, v) \tag{1}$$

In Equation (1), \mathcal{L} denotes the objective function, u denotes the parameters of the generator and v denotes the parameters of the discriminator. In the rest of this paper, G denotes the generator, and D denotes the discriminator. In GAN, x refers to the samples in the real distribution being learned, and z is the target distribution. We set the optimization goal of G to maximize f_G and the optimization goal of D to maximize f_D . In this way, in GAN, we can represent the process as a dynamical system using the ordinary differential equation, and the parameters of this dynamical system consist of two parts: $\theta = (u, v)$, then Equation (1) combined with the basic equation of GAN can be written as:

$$\mathcal{L}(u,v) = E_{(x,z)\sim\mathcal{P}}[\mathcal{L}(u,v;x,z)] = E_{x\sim p_{data}(x)}[f(D_v(x))] + E_{z\sim p_z(z)}[f(1-D_v(G_u(z)))]$$
(2)

In Equation (2), *x* denotes the training samples and *z* denotes the noise vectors. GAN is not searching for a globally optimal solution but a local optimal solution. We want the trajectory of the dynamical system to enter a local convergence point, i.e., Nash equilibrium, with continuous iterations, and define the Nash equilibrium point as: $\overline{u} = \arg \max f_G(u, \overline{v})$,

 $\overline{v} = \underset{v}{\arg \max g_D(\overline{u}, v)}$ It is easy to prove that for a zero-sum game (f = -g), at the Nash equilibrium point, its Jacobian matrix (Equation (3)) is negative definite.

$$\begin{bmatrix} \nabla_u^2 f_G(u,v) & \nabla_{u,v} f_G(u,v) \\ \nabla_{v,u} f_D(u,v) & \nabla_v^2 f_D(u,v) \end{bmatrix}$$
(3)

We can determine whether local convergence is reached by checking the properties of the Jacobian matrix. If at some point, its first-order derivative is 0 (as Equation (4)) and its Jacobian matrix is negative definite, then that point is a Nash equilibrium point.

$$\begin{pmatrix} 0\\0 \end{pmatrix} = \begin{pmatrix} \nabla_u f_G(u,v)\\ \nabla_v f_D(u,v) \end{pmatrix}$$
(4)

We know that the eigenvalues of semi-negative definite matrices are all less than or equal to zero. Then: if the eigenvalues of the Jacobian matrix at a point are all negative real numbers, the training process converges with a sufficiently small learning rate; if the eigenvalues appear to be complex, the training does not achieve local convergence in general; if the real part of the complex eigenvalues is small and the imaginary part is relatively large, some very demanding learning rate is needed to reach convergence. To deal with this situation, we first analyzed the connection and difference between using the Hessian penalty and the Jacobian discriminant (Section 3.2) and approached this convergence with the Hessian–Jacobian consistent orthogonal method, which is partially seen in Section 3.3.

3.2. Hessian Penalty and Jacobian Penalty

The Hessian penalty and the Jacobian penalty each have their own focus in characterizing the original data distribution. This section focuses on the parts of these two things that we use.

3.2.1. Hessian Penalty

Now consider the Hessian matrix, first making the non-diagonal of the Hessian matrix zero, so that its partial derivatives only act independently in each direction, as Equation (5).

$$H_{ij} = \frac{\partial^2 G}{\partial z_i \partial z_j} = \frac{\partial}{\partial z_j} \left(\frac{\partial G}{\partial z_i} \right) = 0$$
(5)

The purpose of this is that z_j has no effect on how the perturbation z_i changes the output of *G*. The Hessian matrix thus corresponds to the loss $\mathcal{L}_H(G) = \sum_{i=1}^{|z|} \sum_{j\neq i}^{|z|} H_{ij}^2$, but this is difficult to solve directly for the calculation, and we need to manipulate it further. After estimating it with unbiased empirical variance, we have Equation (6).

$$\mathcal{L}_H(G) = Var_v \left(v^T H v \right) \tag{6}$$

where v is the Rademacher vector (each entry has the same -1 or +1) and $v^T H v$ is the second-order directional derivative of G in the direction v times |v|. Equation (6) can be estimated using the unbiased empirical variance. In practice, we sample a small number of v vectors (usually only two) to compute this empirical variance. The minimizing Equation (6) is equivalent to minimizing the sum of squares of the non-diagonal elements of the Hessian

matrix. We quickly calculate the second-order directional derivative term in Equation (7) by the second-order central finite difference approximation.

$$v^{T}Hv \approx \frac{1}{\varepsilon^{2}}[G(z+\varepsilon v) - 2G(z) + G(z-\varepsilon v)]$$
(7)

where $\varepsilon > 0$ is a hyperparameter that controls the granularity of the second directional derivative estimate. In our implementation, we use $\varepsilon = 0.1$. Such a Hessian matrix is a square matrix composed of second-order partial derivatives of real-valued functions whose independent variables are vectors. This matrix focuses more on the information change rate of the inverse in the calculation, which corresponds to the extraction and fitting of the surface information (compared to the Jacobi matrix). The Jacobian matrix is a matrix in which the first-order partial derivatives are arranged in a certain way to obtain the Hessian matrix in the shallow features. The Jacobian Penalty corresponding to the Jacobian matrix is described in Section 3.2.2.

3.2.2. Jacobian Penalty

Focusing on the Jacobian matrix, and to more clearly reflect the role of the input parameters and the implied variables on the model, the previous equation $f_G(u, v)$ in Equation (3) is now replaced by $f_G(x, z)$. It is well known that in such a deep generative model, if the correlation coefficients of changes due to different latent dimensions are zero, i.e., $Cov(z_i, ..., z_j) = 0$, then it is considered that each potential dimension controls the variation of only one variation factor. Using the Jacobian matrix mentioned in the kinetic Equation (4), combined with [14], using the Jacobian vector and constraining it to be orthogonal in different dimensions, we can obtain Equation (8).

$$J_{ij} = \left[\frac{\partial f_G(x,z)}{\partial z'_i}\right]^T \frac{f_G(x,z)}{\partial z'_j} = 0$$
(8)

In Equation (8), $\partial z'_i$ and $\partial z'_j$ represents the parameters of the function $f_G(x, z)$ in different directions in the *Z* space. Then the *G* loss of Jacobian penalty becomes the sum of the squared term of Equation (8), and to reduce the computational effort following [43], we rewrite the formula using the Hutchinson estimate as Equation (9).

$$\mathcal{L}_{J}(G) = Var_{v} \Big[(j_{d}v)^{T} j_{d}v \Big]$$
⁽⁹⁾

In Equation (9), we use $j_d = \frac{\partial f_G(x,z)}{\partial z}$ to represent a Jacobian vector. v is the Rademacher vector (each entry has an equal probability of -1 or 1) and *Var* denotes the variance. $j_d v$ is the first-order directional derivative in the v times |v| direction. $j_d v$ can be efficiently computed by the first-order finite difference approximation as Equation (10).

$$\dot{j}_d v = \frac{1}{\varepsilon} [G(z + \varepsilon v) - G(z)]$$
(10)

The operation of this step (Equation (10)) is very similar to the simplification of the Hessian matrix ((Equation (7)), which allows our subsequent calculations to be simplified with similar operations, and it is easy to associate them with each other and the characteristics of their application on GAN networks. We will explain this in Section 3.3.

3.3. Association Between the Hessian Penalty and the Jacobian Penalty

In Section 3.2, we have derived through formulas and calculations the Jacobian penalty (Equation (9)) and the Hessian penalty (Equation (6)). Although they target first-order and second-order information on the solution domain, respectively, they are surprisingly consistent in the solution process, which makes practical calculations convenient. We have also analyzed the association between the two for this case. We give our loss at the end of this subsection.

The relationship between the Hessian matrix and the Jacobian matrix is first given: the Hessian matrix is the inverse of the Jacobian matrix of the function's gradient. The formula expression is in Equation (11).

$$\left(H_f(p)\right)_{ij} = \frac{\partial}{\partial x_j} (\nabla f(p))_i = \frac{\partial^2 f}{\partial x_j \partial x_i} (p) = \left(J_{\nabla f}(p)\right)_{ij} \tag{11}$$

In Equation (11), f is a function from that transforms points into column vectors. p is is the matrix parameters, $H_f(p)$ denotes the hessian matrix, and $J_{\nabla f}(p)$ denotes the Jacobi after the gradient of the matrix f. Thus, the relationship between the Hessian matrix and the Jacobian matrix can be abbreviated as $H(f(x)) = J(\nabla f(x))^T$, considering the nature of matrix operations, so that transposition of the matrix is required in the actual computation.

The Hessian penalty encourages the generator to use the diagonal Hessian of the output concerning the input, using the orthogonalized features as the target of the model, approximated by an unbiased stochastic approximator. The Hessian penalty [13] uses a second-order central finite difference approximation. In contrast, a similar approximation method appears in the same [14], where the Jacobian penalty focuses more on the matrix composed of Jacobian vectors. The orthogonalized features are used as the model's objective to obtain the corresponding Jacobian objective function. This Jacobi approximation calculation used by [14] differs little from the Hessian penalty.

Approximate Computational Procedure and Loss Function

Considering the Taylor expansion of f(x), we note that the Hessian matrix appears in the local gradient expansion, and we add the Hessian as a linear operator to the expansion to obtain Equation (12)

$$\nabla f(x + \Delta x) = \nabla f(x) + H(x)\Delta x + \mathcal{O}\left(\|\Delta x\|^2\right)$$
(12)

where $\nabla f(x)$ is the gradient of f(x), $H(x)\Delta x$ is the f(x) incremental value about Δx , and the tail term $\mathcal{O}(||\Delta x||^2)$ is the second-order infinitesimal term about Δx . To simplify the calculation, we vectorize Δx , so that $\Delta x = \varepsilon v$, substituting into the equation $H(x)\Delta x$, we have Equation (13).

$$H(x)\Delta x = H(x)\varepsilon v = \varepsilon H(x)v = \nabla f(x+\varepsilon v) - \nabla f(x) + \mathcal{O}(\varepsilon^2)$$
(13)

In Equation (13), ε is the scalar and v is the vector, and dividing both sides of Equation (13) by ε , we have $H(x)v = \frac{1}{\varepsilon}[\nabla f(x + \varepsilon v) - \nabla f(x)] + O(\varepsilon)$, and removing the infinitesimal term in Equation (13), the formula is equal to Equation (14).

$$H(x)v \approx \frac{1}{\varepsilon} [\nabla f(x + \varepsilon v) - \nabla f(x)]$$
(14)

The approximation of the removal of the term($\mathcal{O}(\varepsilon)$) usually brings numerical instability, we need to make $\mathcal{O}(\varepsilon)$ small enough, and the effect of $\mathcal{O}(\varepsilon)$ can be ignored in actual calculations, $\mathcal{O}(\varepsilon)$ is very small, so in the actual calculation, we remove this . The effect of Equation (14) and the calculation procedure is the same as that of the original one.

In order to apply the present method to GAN training, we only need to modify the loss of G as Equation (15). Referring to the derivation of Equation (6) for $\mathcal{L}_H(G)$ and Equation (9) for $\mathcal{L}_J(G)$, combined with GAN's original $\mathcal{L}_{G_o} = \underset{z \sim p_z(z)}{\mathbb{E}} [f(1 - D(G(z)))]$, we

can obtain Equation (15). Here, the values of λ_1 and λ_2 are calculated using game theory. For the specific solution process, refer to Appendix A.1. Likewise, the two parameters of this part are also used in the network architecture; for the network architecture algorithm pseudo-code, refer to Appendix A.2.

$$\mathcal{L}_{G} = \mathop{\mathrm{E}}_{z \sim p_{z}(z)} [f(1 - D(G(z)))] + \lambda_{1} \mathop{\mathrm{E}}_{z \sim p_{z}(z)} [\mathcal{L}_{H}(G)] + \lambda_{2} \mathop{\mathrm{E}}_{z \sim p_{z}(z)} [\mathcal{L}_{J}(G)]$$
(15)

In Equation (15), λ_1 and λ_2 are hyperparameters. The inclusion of $\mathcal{L}_H(G)$ and $\mathcal{L}_J(G)$ in GAN training is beneficial to learn the representation and to achieve the generation of controllable images. The specific calculations for the hyperparameters λ_1 and λ_2 are shown in Appendix A.1. We found that using these hyperparameters only to adjust the training parameters in the model architecture, without using this disentanglement loss in Equation (15), can mitigate the pattern collapse. The use of the hyperparameters in the model is referenced in Appendix A.2.

3.4. Network Model Structure

Figure 1 shows the network model proposed in this paper, where the green solid line frame is the original model architecture, and the red dotted lines are the parts added by our method. The green part divides the generative adversarial network training into two processes. The first process updates the generation loss, and the second process updates the discriminant loss. GLoss and DLoss correspond to the loss functions of the generator and discriminator, respectively. Our method is divided into two parts to be added to the model; one is the Update red dashed box in process one and process two, which is mainly used to mitigate mode collapse; the other is the red dashed box corresponding to Gloss. We use Hessian combined with Jacobian for disentangled representation. The two added parts of our model are independent of each other. The red dashed part in the green box is used alone for mode collapse mitigation, and the red dashed part out of the green box is used alone for disentanglement representation.



Figure 1. Structure of the network model.

3.5. Evaluation Metrics

Inception Score (IS) [44] uses an image class classifier to evaluate the quality of the generated images. The image class classifier used is Inception Net-V3. IS is a measure of the clarity and diversity of the generated images, and the bigger the IS value, the better. The specific formula is as Equation (16).

$$\mathrm{IS}(G) = \exp\left(\mathbb{E}_{x \sim p_g} D_{KL}(p(y \mid x) \| p(y))\right)$$
(16)

In Equation (16), D_{KL} denotes KL divergence (Kullback–Leibler divergence). For a picture *x*, the probability distribution that belongs to all classes is p(y | x). p(y) denotes marginal probability.

Fréchet Inception Distance (FID) [30] does not rely on a classifier, but directly considers the distance between the generated data and the real data at the feature level. FID uses the 2048-dimensional vector before the full connection of Inception Net-V3 as the feature of the

picture. FID is the distance between the generated image and the real image; the smaller the FID value, the better. The specific formula is as Equation (17).

$$FID = \left\|\mu_r - \mu_g\right\|^2 + Tr\left(\Sigma_r + \Sigma_g - 2\left(\Sigma_r \Sigma_g\right)^{1/2}\right)$$
(17)

In Equation (17), μ_r denotes the feature mean of the real images. μ_g denotes the feature means of the generated images. Σ_r denotes the covariance matrix of the real images. Σ_g denotes the covariance matrix of the generated images. Tr denotes the trace of the matrix.

Perceptual Path Length (PPL) [1] subdivides the interpolation path of two noise points into multiple segments, finds the length of each segment, and then averages it. PPL evaluates the distance at which the generator changes from one image to another. The smaller the PPL value, the better. The specific formula is as Equation (18).

$$PPL = \mathbb{E}\left[\frac{1}{\varepsilon^2}d(G(\operatorname{slerp}(z_1, z_2; t)), G(\operatorname{slerp}(z_1, z_2; t + \varepsilon)))\right]$$
(18)

In Equation (18), ε denotes subdivision, replaced with 1×10^{-4} . *d* denotes perceptual distance, measured using pre-trained VGG. *G* denotes the image generator. Function slerp denotes spherical linear interpolation, an interpolation method. *t* denotes the interpolation parameters, subject to the uniform distribution.

4. Experiments

We conducted experiments on datasets such as MNIST, Cifar-10, and CIEVR with several different generative adversarial network models. We evaluate generation based on two criteria: model collapse and generated sample quality. Model collapse and generated sample quality are the two criteria we target and evaluate by Fréchet Inception Distance (FID) [30] and Inception Score (IS) [44]. An important metric for assessing model performance is IS. Calculating IS requires Inception Net-V3 (the third version of Inception Net), and the FID calculates the real samples and generates the distance between samples in the feature space. Different from IS, a lower FID implies higher image quality and diversity.

A comparative test of GANs and the original method is performed to give experimental curves and correlation analysis. Experiments with λ_1 and λ_2 were selected, and then our experiments were repeated on other GANs with quantitative analysis. The same architecture and hyperparameters for each method separately ensure fairness in the experiments.

In the disentanglement part, we set all the image sizes to 128×128 according to the training arrangement in Hessian. The tested datasets are Edge + Shoes and CLEVR datasets, respectively.

In Section 4.1, we introduce the datasets used for the experiments, and in Sections 4.2 and 4.3, we test our approach in model collapse and disentanglement.

4.1. Introduction to the Datasets

The MNIST dataset was sourced from a mixture of two datasets, one handwritten by employees from the Census Bureau and one handwritten by high school students; this dataset had 60,000 training samples and the rest 10,000 were test samples, with each image being a single number from 0 to 9. Example images of the dataset is shown in Figure 2a.

The Cifar-10 dataset has 60,000 images, which are 32×32 and divided into ten categories with 6000 images in each category. Example images of the dataset is shown in Figure 2f.

CelebA is the open dataset of the Chinese University of Hong Kong, containing 202,599 images of 10,177 celebrity identities. Each image is annotated with 40 attributes. Example images of the CelebA dataset are shown in Figure 2d. CelebA-HQ is a high-quality face image of CelebA, consisting of 30,000 images. In the experiments on the CelebA-HQ dataset, we mainly use 256×256 size images. Example images of the CelebA-HQ dataset are shown in Figure 2e.

The CLEVR dataset is known as Compositional Language and Elementary Visual Reasoning. CLEVR contains 100,000 rendered images and approximately 1 million automatically generated questions, of which 853,000 are mutually exclusive. Example images of the dataset are shown in Figure 2c.

Edge + Shoes is a simple dataset. This is a large dataset that includes 50,025 images. These images are divided into four categories: shoes, sandals, flip-flops, and boots. Example images of the dataset are shown in Figure 2b.



Figure 2. The datasets used in this article.

4.2. Resistance Model Collapse

In this part, we do not directly use the loss of the Formula (15), but instead use the λ_1 and λ_2 parameters in the loss, which are used in the model, specifically using the part of the algorithm presented in Appendix A.2. The specific use of λ_1 and λ_2 in the model is in Algorithm A1. The reason for doing so is to optimize the current GAN model and make the basis for the disentanglement.

We run real image generation experiments on different datasets: MNIST, CIFAR-10, and CelebA. It has been demonstrated experimentally that our approach is general to some generative models.

We train all models for 25,000 iterations, and in each iteration, we pass the parameters in the generator to the discriminator at a certain frequency. At inference, we generate 10k samples from each training model and measure these two metrics. We report in Table 1 the average scores (IS) of the images generated on the MNIST dataset [34], and p means *p*-value of the statistical significance. As shown in Table 1, ours outperforms the other methods. We divided the original training datasets into two parts to measure their IS. Since the IS values calculated by different sample segmentation methods are different (this is because the calculation of IS needs to calculate the two groups of pictures, and the segmentation method determines the difference of the calculated values), so it is not given in the Tables 1 and 2. The first *p*-value is determined from the IS means of the original method and the IS means within the original training datasets. The second *p*-value (*p* (ours)) is determined by the IS means of our method and the IS means of the original training datasets. These two *p*-values compare the IS means of the method before and after the improvement with the original training datasets. Taking $\alpha = 0.001$ as the judging criterion, our results show that the difference between the IS means of the images generated by our GANs network and the IS means of the original datasets are statistically significant.

GAN	IS (Mean)	IS (Std)	р	IS (Ours Mean)	IS (Ours Std)	p (Ours)
BEGAN [28]	1.366	0.179	< 0.001	1.388	0.242	< 0.001
BGAN [36]	1.250	0.175	< 0.001	1.372	0.187	< 0.001
EBGAN [10]	1.356	0.194	< 0.001	1.371	0.241	< 0.001
INFOGAN [32]	2.054	0.171	< 0.001	2.070	0.207	< 0.001
CGAN [31]	2.022	0.196	< 0.001	2.302	0.291	< 0.001
AAE [15]	2.289	0.298	< 0.001	1.995	0.251	< 0.001
ACGAN [16]	1.956	0.146	< 0.001	2.041	0.271	< 0.001
LSGAN [9]	1.404	0.183	< 0.001	1.363	0.139	< 0.001
SGAN [17]	1.418	0.159	< 0.001	1.459	0.192	< 0.001

Table 1. Score (IS) test on GANs (MNIST).

The experimental results in Figure 3 show that the BGAN [36] model modified by our method has a Discriminator closer to 0.5 than the original BGAN [36], and the fluctuation is much smaller than that of the model without our method. At the same time, it has faster convergence and stability.



Figure 3. The results of the CGAN [31] and BGAN [36] model.

Figure 3 shows that our improved model is more stable than the original one in CGAN [31]. For more experimental results of the loss function curves refer to Appendix B.

After training all models for 100K iterations, we evaluated the method on CIFAR-10. Unlike MNIST, the models in this dataset are difficult to handle. We need to evaluate color images, and the contents of pictures are more diverse. That is why we use two different metrics (IS and FID) for inferring the generator quality and diversity. As shown in Table 2, our method consistently outperforms all other methods on these two metrics.

We trained on the standard CIFAR-10 dataset and used the Inception v3 image classification model to calculate the similarity of statistical aspects of computer vision features to measure the similarity of the two datasets. As shown in Table 3, our method generates FID values with the optimal ones.

GAN	IS (Mean)	IS (Std)	p	IS (Ours Mean)	IS (Ours Std)	p (Ours)
BEGAN [28]	1.735	0.215	< 0.001	2.005	0.236	< 0.001
EBGAN [10]	1.666	0.224	< 0.001	1.733	0.244	< 0.001
INFOGAN [32]	1.563	0.197	< 0.001	1.677	0.208	< 0.001
BGAN [36]	1.822	0.203	< 0.001	2.932	0.221	< 0.001
CGAN [31]	1.978	0.209	< 0.001	2.065	0.235	< 0.001
AAE [15]	2.012	0.214	< 0.001	1.917	0.199	< 0.001
ACGAN [16]	1.771	0.231	< 0.001	1.871	0238	< 0.001
LSGAN [9]	1.846	0.227	< 0.001	1.959	0.236	< 0.001
SGAN [17]	1.668	0.158	< 0.001	2.005	0.183	< 0.001

Table 2. Score (IS) test on GANs (Cifar-10).

GAN	FID (Original)	FID (Ours)
BGAN [36]	243.633	226.164
BEGAN [28]	417.625	377.760
CGAN [31]	382.941	373.581
AAE [15]	187.612	181.169
ACGAN [16]	231.143	249.174
LSGAN [9]	149.675	133.783
SGAN [17]	147.770	130.251

Table 3. Score (FID) test on GANs (Cifar-10).

In Figure 4, we compare the image generated in CelebA by the original DCGAN [33] and ours. We reduced the image size of CelebA to 96×96 to speed up the training process for faster results. The generated image is also in the size of 96×96 . Compare our method with the method before improvement in Figure 4, it is not difficult to find that the image face generated by the original DCGAN [33] is blurry and distorted, and the hairstyle has fewer changes in shape and eyebrow shape. Figure 4b is generated by our method, which is improved on the original basis, it has better clarity and diversity. In Figure 4, the faces of the upper character (with a variety of hairstyles) in the (b) pictures are more clear, and the forehead hair of the task in the lower (b) pictures are from long to short, showing the diversity of hairstyles. The image generated by our method also has fewer facial artifacts than the images on the original DCGAN (a).



Figure 4. Graph generated without our Network(DCGAN [33]) (a), with ours (b).

4.3. Disentanglement Experiments

4.3.1. Qualitative Experiments

In this subsection, we compare Hessian penalty [13] and OroJaR [14] with our method. After adding the λ_1 and λ_2 parameters in the model training, some cases will bring more serious attribute entanglement, so we add the Equation (15) to disentanglement and compare the current best results.

Figure 5 shows experiments been done on fox, balloons, and Tibetan mastiffs on ImageNet. We moved the images in Z-space for each row. Z-space represents the sample space of latent variables. The input position of the Z-space refers to Figure 1. Through the change of the Z-space parameters, the model generates a series of pictures, as shown in Figure 5. The role of these diagrams is explained next. The first line in Figure 5 shows the enlargement of the fox's face, after disentanglement, the Z-space parameters can control the distance between the object and the lens; the second line shows the separation and change of the background, we separate the balloon from the background, separate the balloon property, and change the background without changing this property, and the third line shows the separation and action of the foreground object (the Tibetan mastiff gradually stands up). The foreground object is a Tibetan mastiff, and the background is grass. By selecting the direction of the Z-space, the foreground object can make actions without changing the background.



Figure 5. The orthogonal orientation learned on BigGAN [2].

In Figure 6, we present the experimental results of our model on the Edge + Shoes dataset. We choose the two interpretable dimensions of shoe color and style to display, with each row corresponding to one interpretable dimension. Figure 6 shows OroJaR [14], Hessian penalty [13], and our method learned the two separated changes, i.e., style and shape of shoes (the first row of each sub-image is generated by the Hessian method [13], the middle row by the OroJaR [14] method, and the last row is produced by our method). Subfigure (a) shows the performance of different methods in disentangling shape entanglement, and our method covers more shapes. Subfigure (b) shows different methods in disentangling color in terms of entanglement performance; Hessian penalizes [13] to generate a single color of the image, and OroJaR [14] generates a serious color coupling of shoe images. Our method covers more colors, and the independence of colors is better than others.



Figure 6. Comparison of the disentanglement quality of OroJaR [14], Hessian penalty [13] and our method for Edge + Shoes.

Figure 7 shows the results on the CLEVR simple dataset. Our method is able to separate the potential space effectively. We show the four scoring dimensions of *x* axial shift, *y* axial shift, color, and shape. Hessian penalty learning independently controls the vertical and horizontal axial position of the object (first row of each subplot of Figure 7). However, the shape is not controlled during the vertical movement. OroJaR [14] successfully separates the horizontal, shape factors in CLEVR (second row of each subfigure of Figure 7), but shows a tendency to shift toward the horizontal in the separation of the vertical movement. Lastly, our method can successfully separate independent axial shifts and colors in CLEVR (the third row of each subplot of Figure 7).



Figure 7. Comparison with OroJaR [14], Hessian penalty [13], and our method for the disentanglement quality of CLEVR Sample.

Figure 8 shows the comparison on the CLEVR complex dataset. Here we show the relative position as a factor. From Figure 8, we can conclude: (1) The Hessian penalty does not perform well in the relative positions of two objects, and when the Hessian penalty changes the relative positions of two objects, the shapes or colors of these two objects change at the same time (first row of each subplot of Figure 8). (2) OroJaR [14] changes its shape when the relative position changes (second row of each subfigure of Figure 8). (3) Our method does not change shape and color when the relative position changes (third row of each subfigure of Figure 8).



(a) Change relative position

(b) Change size

Figure 8. Comparison of disentanglement the quality of the CLEVR complex by our method, Hessian penalty [13], and OroJaR [14].

In Figure 9, subfigure (a) shows that, in the case of adding objects, it tries to keep the color properties unchanged. Subfigure (b) shows that it is possible to add objects and keep their shape properties in the case of two colors. The graphs in the first row of subgraphs (a) and (b) in Figure 9 are generated by Hessian penalty [13], the graphs in the second row are generated by OroJaR [14], and the graphs in the third row are generated

by our method. The graphs generated by the Hessian penalty [13] appear in five colors as the objects gradually increase from one to six. The graphs generated by OroJaR [14] appear in four colors. Our method emerges in two colors. Although none of the three methods achieves the best-case scenario of producing only one color, our method is more advantageous in color preservation when adding objects. In subgraph (b), the graphs generated by Hessian penalty [13] appear in two colors and two shapes, and the graphs generated by OroJaR [14] appear in more than two colors. However, the shape remains relatively stable (five cylinders in a row) compared to the previous two methods. The shape-preserving effect of our method is also stable in two colors.



Figure 9. Comparison of disentanglement the quality of the CLEVR dataset by our method, Hessian penalty [13] and OroJaR [14] for adding objects.

We conduct disentanglement experiments on the CelebA-HQ dataset and compare the performance of Hessian penalty [13], OroJaR [14], and our method on face rotation. In Figure 10, the graphs in the first row are generated by Hessian penalty [13], the graphs in the second row are generated by OroJaR [14], and the graphs in the third row are generated by our method. The properties required for disentanglement are independent of each other. When one property is changed, the other properties remain unchanged. From the comparison in Figure 10, it can be found that, when the Hessian penalty [13] modifies the rotation attribute, the eyes of the characters in the figures always look towards the camera and do not rotate with the characters, resulting in disentanglement. Moreover, artifacts are created at the edges in the first row of images. The eyes of the OroJaR [14] method in the second line rotate with the picture character, but when only controlling the rotation of the character's avatar, the mouth shape of the second line changes uncontrollably (from normal to slightly flared). The third row shows the performance of our method on face



rotation. Our method achieves the rotation of the eyes and the character together, and the mouth shape remains unchanged, which cannot be achieved by the first two methods.

Figure 10. Comparison of rotation disentanglement of the CelebA-HQ dataset by our method, Hessian penalty [13], and OroJaR [14].

4.3.2. Quantitative Evaluation

In this subsection, we compare our method quantitatively with some models. Referring to [13], we use perceptual path length (PPL) [1], Frechet inception distance (FID) [30], and variance predictability disentanglement measure (VP) [45] as quantitative metrics. Lower FID and PPL are better, and higher VP is better.

As can be seen from Table 4, our method has obtained some better results on some datasets, which indicates that our method has a better disentanglement representation.

	Ec	Edges + Shoes		CLEVR-Simple		CLEVR-Complex			CLEVR-U			
Method	PPL (↓)	FID (↓)	VP (†)	PPL	FID	VP	PPL	FID	VP	PPL	FID	VP
InfoGAN [32]	2952.2	10.4	15.6	56.2	2.9	28.7	83.9	4.2	27.9	766.7	3.6	40.1
ProGAN [46]	3154.1	10.8	15.5	64.5	3.8	27.2	84.4	5.5	25.5	697.7	3.4	40.2
SeFa [12]	3154.1	10.8	24.1	64.5	3.8	58.4	84.4	5.5	30.9	697.7	3.4	42.0
Hessian Penalty [13]	554.1	17.3	28.6	39.7	6.1	71.3	74.7	7.1	42.9	61.6	26.8	79.2
OroJaR [14]	236.7	16.1	32.3	6.7	4.9	76.9	10.4	10.7	48.8	40.9	4.6	90.7
Ours	336.79	12.5	33.7	27.6	3.3	79.1	36.3	4.3	49.3	44.1	3.3	88.3

Table 4. Comparison of different methods on Edge + Shoes and CLEVR.

4.4. Limitation and Discussion

Our proposed method performs well on mode collapse mitigation and disentanglement, scoring higher than other methods on the test dataset. However, our mode collapse mitigation and disentanglement methods need to be used separately and cannot be combined with one model simultaneously. Our approach to mode collapse mitigation focuses on making the discriminator and generator of the generative adversarial network interact with weights, and discriminator and generator can achieve better results in this way. More features will also bring the coupling of properties, making it impossible to achieve the expected disentanglement effect after adding the disentanglement loss. In the future, applying both mode collapse mitigation and disentanglement methods to the model is a considered direction for improvement.

5. Conclusions

In this work, we introduced a new method to train generative networks. We apply our method to generative adversarial training by learning disentanglement representations. We train the generator to optimize this disentanglement representation through loss or hyperparameters. Our framework accumulates many desirable properties: it does not require any additional trainable parameters, it operates in an unsupervised environment, and it outperforms state-of-the-art methods in some image datasets. Moreover, our method exhibits stable adversarial training, and our method allows embedding any variant of the GAN model. In the future, the neural radiance field can be combined with existing models to achieve a better disentanglement representation.

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Abbreviations

The following abbreviations are used in this manuscript:

- GAN Generative Adversarial Networks
- D Discriminator
- G Generator
- KL Kullback-Leibler divergence
- FID Fréchet Inception Distance
- IS Inception Score

Appendix A. Supplementary Theory

Appendix A.1. Eagle Dove Game

Consider the later stage of the training process, when the Generation distribution of G is almost the same as the original distribution, the corresponding equilibrium will be similar. At this time, DG and the corresponding equilibrium are likely interlaced and matched, so only this is difficult for such a theory to describe the DG game process completely. Thus, this model was explained from the perspective of game theory. "Eagle Dove Game" is a classic game model. The basic description of this model is that, in a particular game environment, both sides of the game have the right to be an eagle and a dove. If both sides of the game choose eagles, then they go out to discount the benefits, and the remaining benefits are divided equally. If one party chooses the dove and the other side still chooses the eagle, the other side will win. If both sides choose the doves, compared to the eagle, they could evenly split the gains without losing the competitive discount. Then, the resulting payout matrix is as Table A1:

Table A1. A result payout matrix of Eagle Dove Game.

Strategy	Eagle	Dove
Eagle	S = (v - c)/2, S = (v - c)/2	T = v, P = 0
Dove	P = 0, T = v	R = v/2, R = v/2

Table A1 shows the conclusion that the abilities of both parties in the game are the same, v is the benefit and v is the conflict cost paid by both parties, S is the benefit of both parties when both parties choose eagle strategy, T is the benefit of both parties when both parties choose eagle dove strategy, P is the benefit of both parties when both parties choose eagle dove strategy, and *R* is the benefit of both parties when both parties choose eagle strategy. If the skills of both parties are the same, then the choice is good, but just like the relationship between the Discriminator and the Generator, the ability between the two parties in the game is a big gap. Suppose the skills of the two parties differ significantly, in that case, the strategy chosen by the two parties, the Discriminator and Generator must be affected by the ratio of the capabilities of the two parties. When the two parties fight at this time, the obtained gain is related to the difference in the strength of the two parties. It proves that the better Generators trained initially and the Discriminators with average learning speed cannot effectively train the ideal results. Let the ability ratio of the two parties involved in the game to the degree of asymmetry be k: (1 - k), then if the conflict loss is greater than the cooperation gain, the payment matrix is shown in Table A1 can be converted to Table A2:

Table A2. A new result payout matrix of Eagle Dove Game.

Strategy	Eagle	Dove
Eagle	(v-c)/4k, (v-c)/4k	(v, 0)
Dove	(0, v)	(kv, (1 - k)v)

The reason for multiplying by 1/4 in Table A2 is that, when *k* is 0.5, the model degenerates to a classic eagle–dove game. Under the change of payment matrix, neither party in the game has a purely strict dominant strategy equilibrium. Assuming that the Generator's frequency adopting the dove strategy is *x*, and the Discriminator adopts the dove strategy as *y*, then correspondingly, the frequency of the Generator adopts the eagle strategy is 1 - x, and Discriminator adopts the dove strategy as expected 1 - y. Then, according to Table A2, it shows that the expected return of the Generator choosing the eagle strategy is: $\overline{U_A}(x, y) = (1 - x)U_A^H + xU_A^D$ The expected return of the Generator's choice of dove strategy as: $\overline{U_B}(x, y) = (1 - y)U_B^H + yU_B^D$. The Discriminator chooses the expected return of the eagle strategy as: $\overline{U_B}(x, y) = (1 - x)U_B^H + (1 - x)\frac{v-c}{4(1-k)} + xv$. The expected return of the Discriminator chooses the expected return of the corresponding Generator is: $U_A^H = (1 - y)\frac{v-c}{4k} + yv$. In the same way, the average expected return of the corresponding Discriminator is: $U_A^D = (1 - y) \cdot x + ykv$. Let $x = \lambda_1$, $y = \lambda_2$, then, the Nash equilibrium can be solved as:

$$\left((\lambda_1, 1 - \lambda_1), (\lambda_2, 1 - \lambda_2)\right) = \left(\left(\frac{c - v}{M}, 1 - \frac{c - v}{M}\right)\left(\frac{c - v}{M}, 1 - \frac{c - v}{M}\right)\right)$$
(A1)

 $M = c - v + 4kv - 4k^2v$. In this way, we obtain the parameters about DG: λ_1 , λ_2 . These parameters can provide us with important hyperparameters initialization guidance in subsequent actual experiments.

Appendix A.2. Model Architecture

In our model, the DG process is regarded as two parts. One part is to update the Generator's weights; the other part is to update the Discriminator's weights. The direction of the Generator and the Discriminator on the saddle point is opposite, the point that finally meets is the saddle point, which is also the Nash equilibrium point of the game. Updating the weights of the Generator and the Discriminator will cause their updated weight direction to be opposite the original direction simultaneously. Its pseudo-code refers to Algorithm A1.

Require:	_
Images in datasets:	
k, λ_1 , λ_2 ; k is a default GAN training steps, λ_1 is D interaction frequency, λ_2 is	D
interaction frequency;	
Ensure:	
Distribution of target datasets;	
for training iterations do	
for k steps do	
Take samples from noise prior.	
Take examples from data generating distribution.	
Update the D.	
for $k\lambda_1$ steps do	
Update the G.	
end for	
Take examples from data generating distribution.	
for $k\lambda_2$ steps do	
Update the G.	
end for	
Take examples from data generating distribution.	
end for	
Take noise samples from noise prior.	
for $k\lambda_1$ steps do	
Take examples from data generating distribution.	
end for	
Update the G.	
end for	

Appendix B. Other Experiments

Both the discriminator and generator of DCGAN [33] use a convolutional neural network to replace the multi-layer perceptron in GAN, and use convolution and deconvolution to replace the pooling layer. The output layer of the generator uses the Tanh activation function, and the other layers use RELU, the discriminator LeakyReLU activation function is used for all layers of. After the network is added to our method, Figure A1 shows that the generator fluctuation of the network is significantly reduced, and the numerical distance between the network and the discriminator is also getting closer.



Figure A1. The results of the DCGAN [33] and BEGAN [28] model.

In Figure A1, we found that our method is not very effective for the network of the GAN network generator that uses the loss of Wasserstein distance, but in the actual training process, it can reach the training completion standard at a faster speed, so that we need early stop to finish training.

A pull-away term was proposed to prevent Generator from generating the same image. In Figure A2, our method has smaller oscillation and convergence on this EBGAN [10], and the time to reach stable generation is also shorter.

INFOGAN [32] divides Z into fixed distribution noise and c some hidden variable information on the basis of DC-GAN. In Figure A2, after adding our method, DG converges faster and gets closer.

The method of LSGAN [9] is to change the objective function of GAN from cross entropy loss to least square loss to improve the image quality and stabilize the training process. After our method is added in Figure A3, the improvement on this GAN is not obvious.



Figure A2. The results of the EBGAN [10] and INFOGAN [32] model.



Figure A3. The results of the LSGAN [9] and SGAN [17] model.

SGAN [17] extends the Generative Adversarial Network (GAN) to semi-supervised context by forcing the discriminator network to output class labels. From Figure A3, we can conclude that our generator is still closer to 0.5.

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