- Convexified CNNs to removing the weight nonlinearity.
- Introduction to Generative Adversarial Networks
  - Inverse conv net for generating images
  - The adversarial game
  - Applications and improved training strategies, WGANs

## Convexifying the parameters pt. 1 (Zhang et al. $16'^1$ )

Consider a two layer convolutional neural network composed of one convolutional layer followed by a fully connected layer. Rather than working with x directly, form P vectors  $z_p(x)$  for p = 1, ..., P where  $z_p(x)$  is the portion of x on patch p of the convolutional layer. Then the  $k^{th}$  component of  $H(x, \theta)$  is given by

$$H(x,\theta)_k = \sum_{j=1}^r \sum_{p=1}^p \alpha_{k,j,p} \sigma(w_j^T z_p(x)).$$

Alternatively if we exclude the nonlinearity we can express this sum by

$$\sum_{j=1}^{r} \sum_{p=1}^{p} \alpha_{k,j,p} \sigma(w_j^T z_p(x)) = \sum_{j=1}^{r} Z(x) w_j$$

where Z(x) has  $z_p(x)$  as its  $p^{th}$  row.

<sup>1</sup>https://arxiv.org/pdf/1609.01000.pdf

## Convexifying the parameters pt. 2 (Zhang et al. $16'^2$ )

Using the trace formula this can be further condensed to

$$H(x,\theta)_k = \operatorname{tr}\left(Z(x)\left(\sum_{j=1}^r w_j \alpha_{k,j}^T\right)\right) = \operatorname{tr}\left(Z(x)A_k\right)$$



The network parameters are given by  $A_k$  nonlinearity is imposed by the  $A_k$  having rank r, and we can express all of the parameters of the matrix by A which is similarly rank r. <sup>2</sup>https://arxiv.org/pdf/1609.01000.pdf

One can impose the network structure through A, but remove the non-convex rank constraint by replacing a convexification, that is the sum of the singular values of A (Schatten-1, or nuclear, norm).

If the convolutional filters and fully connected rows are uniformly bounded in  $\ell^2$  by  $B_1$  and  $B_2$  respectively, then one can replace then the sum of the singular values of A are bounded by  $B_1B_2r\sqrt{n}$ where n is the network output dimension and the network parameters can be considered by varying the nuclear norm bound between 0 and  $B_1B_2r\sqrt{n}$ .

The resulting learning programme is fully convex and can be efficiently solved. The above can be extended to nonlinear activations and multiple layers, learning one layer at a time.

<sup>&</sup>lt;sup>3</sup>https://arxiv.org/pdf/1609.01000.pdf

|                  | basic | rand   | rot    | img    | img+rot |
|------------------|-------|--------|--------|--------|---------|
| $SVM_{rbf}$ [44] | 3.03% | 14.58% | 11.11% | 22.61% | 55.18%  |
| NN-1 [44]        | 4.69% | 20.04% | 18.11% | 27.41% | 62.16%  |
| CNN-1 (ReLU)     | 3.37% | 9.83%  | 18.84% | 14.23% | 45.96%  |
| CCNN-1           | 2.38% | 7.45%  | 13.39% | 10.40% | 42.28%  |
| TIRBM [38]       | -     | -      | 4.20%  | -      | 35.50%  |
| SDAE-3 [44]      | 2.84% | 10.30% | 9.53%  | 16.68% | 43.76%  |
| ScatNet-2 [8]    | 1.27% | 12.30% | 7.48%  | 18.40% | 50.48%  |
| PCANet-2 [9]     | 1.06% | 6.19%  | 7.37%  | 10.95% | 35.48%  |
| CNN-2 (ReLU)     | 2.11% | 5.64%  | 8.27%  | 10.17% | 32.43%  |
| CNN-2 (Quad)     | 1.75% | 5.30%  | 8.83%  | 11.60% | 36.90%  |
| CCNN-2           | 1.38% | 4.32%  | 6.98%  | 7.46%  | 30.23%  |

Table 1: Classification error on the basic MNIST and its four variations. The best performance within each block is bolded. The tag "ReLU" and "Quad" means ReLU activation and quadratic activation, respectively.

<sup>4</sup>https://arxiv.org/pdf/1609.01000.pdf

## Convexified CNN: CIFAR10 (Zhang et al. 16'<sup>5</sup>)



Table 3: Classification error on the CIFAR- Figure 4: The convergence of CNN-3 and 10 dataset. The best performance within each block is bolded.

CCNN-3 on the CIFAB-10 dataset.

|             | CNN-1  | CNN-2  | CNN-3  |
|-------------|--------|--------|--------|
| Original    | 34.14% | 24.98% | 21.48% |
| Convexified | 23.62% | 21.88% | 18.18% |

Table 4: Comparing the original CNN and the one whose top convolution layer is convexified by CCNN. The classification errors are reported on CIFAR-10.

<sup>5</sup>https://arxiv.org/pdf/1609.01000.pdf

## CNN model through sparse coding (Papyan et al. 16'<sup>6</sup>)

Consider a deep conv. net composed of two convolutional layers:



The forward map (note notation using transpose of  $W^{(i)}$ ):

$$Z_{2} = \sigma \left( b^{(2)} + (W^{(2)})^{T} \sigma \left( b^{(1)} + (W^{(1)})^{T} x \right) \right)$$

<sup>6</sup>https://arxiv.org/pdf/1607.08194.pdf

#### Deconvolutional NN data model (Papyan et al. 16'<sup>7</sup>)



Two layer deconvolutional data model with weight matrices fixed,  $W^{(i)} = D_i$ , and  $\Gamma_i \ge 0$  whose values compose data element X.

<sup>&</sup>lt;sup>7</sup>https://arxiv.org/pdf/1607.08194.pdf

## Generative deep nets (Goodfellow et al. 14'<sup>9</sup>)





Figure 1: DCGAN generator used for LSUN scene modeling. A 100 dimensional uniform distribution Z is projected to a small spatial extent convolutional representation with many feature maps.

<sup>8</sup>https://arxiv.org/pdf/1511.06434.pdf
<sup>9</sup>https://arxiv.org/pdf/1406.2661.pdf

#### Generative deep nets (Goodfellow et al. 14'<sup>10</sup>)

Train the two network parameters using the objective

$$\min_{G} \max_{D} n^{-1} \sum_{\mu=1}^{n} \log(D(x_{\mu}, y_{\mu})) + p^{-1} \sum_{p} \log(1 - D(G(z_{p}), y_{p}))$$

**Algorithm 1** Minibatch stochastic gradient descent training of generative adversarial nets. The number of steps to apply to the discriminator, k, is a hyperparameter. We used k = 1, the least expensive option, in our experiments.

for number of training iterations do

for k steps do

- Sample minibatch of m noise samples  $\{z^{(1)}, \ldots, z^{(m)}\}$  from noise prior  $p_g(z)$ .
- Sample minibatch of m examples  $\{x^{(1)}, \ldots, x^{(m)}\}$  from data generating distribution  $p_{\text{data}}(x)$ .
- Update the discriminator by ascending its stochastic gradient:

$$\nabla_{\theta_d} \frac{1}{m} \sum_{i=1}^m \left[ \log D\left( \boldsymbol{x}^{(i)} \right) + \log \left( 1 - D\left( G\left( \boldsymbol{z}^{(i)} \right) \right) \right) \right].$$

end for

- Sample minibatch of m noise samples  $\{z^{(1)}, \ldots, z^{(m)}\}$  from noise prior  $p_a(z)$ .
- Update the generator by descending its stochastic gradient:

$$\nabla_{\theta_g} \frac{1}{m} \sum_{i=1}^m \log\left(1 - D\left(G\left(\boldsymbol{z}^{(i)}\right)\right)\right).$$

#### end for

The gradient-based updates can use any standard gradient-based learning rule. We used momentum in our experiments.

#### <sup>10</sup>https://arxiv.org/pdf/1406.2661.pdf

#### Generative deep nets (Radford et al. $16'^{11}$ )



Figure 2: Generated bedrooms after one training pass through the dataset. Theoretically, the model could learn to memorize training examples, but this is experimentally unlikely as we train with a small learning rate and minibatch SGD. We are aware of no prior empirical evidence demonstrating memorization with SGD and a small learning rate.

<sup>11</sup>https://arxiv.org/pdf/1511.06434.pdf

# Generative deep nets (Radford et al. 16'12)



Figure 3: Generated bedrooms after five epochs of training. There appears to be evidence of visual under-fitting via repeated noise textures across multiple samples such as the base boards of some of the beds.

<sup>12</sup>https://arxiv.org/pdf/1511.06434.pdf

#### Wasserstein GAN (Arjovsky et al. 17'14)

One of the central challenges with GANs is the ability to train the parameters. Improvements have been made through choice of generative architecture (DC-GAN of Radford) and through different training objective functions (W-GAN)

Algorithm 1 WGAN with gradient penalty. We use default values of  $\lambda = 10$ ,  $n_{\text{critic}} = 5$ ,  $\alpha = 0.0001$ ,  $\beta_1 = 0$ ,  $\beta_2 = 0.9$ .

**Require:** The gradient penalty coefficient  $\lambda$ , the number of critic iterations per generator iteration  $n_{\text{critic}}$ , the batch size m, Adam hyperparameters  $\alpha$ ,  $\beta_1$ ,  $\beta_2$ .

**Require:** initial critic parameters  $w_0$ , initial generator parameters  $\theta_0$ .

1: while  $\theta$  has not converged do

```
2:
               for t = 1, \dots, n_{\text{critic}} do
 3:
                       for i = 1, ..., m do
                               Sample real data \boldsymbol{x} \sim \mathbb{P}_r, latent variable \boldsymbol{z} \sim p(\boldsymbol{z}), a random number \epsilon \sim U[0, 1].
 4:
 5:
                              \tilde{x} \leftarrow G_{\theta}(z)
                              \hat{\boldsymbol{x}} \leftarrow \epsilon \boldsymbol{x} + (1-\epsilon)\tilde{\boldsymbol{x}}
 6:
                               L^{(i)} \leftarrow D_w(\tilde{\boldsymbol{x}}) - D_w(\boldsymbol{x}) + \lambda (\|\nabla_{\hat{\boldsymbol{x}}} D_w(\hat{\boldsymbol{x}})\|_2 - 1)^2
 7:
                       end for
 8:
 9:
                       w \leftarrow \operatorname{Adam}(\nabla_w \frac{1}{m} \sum_{i=1}^m L^{(i)}, w, \alpha, \beta_1, \beta_2)
10 \cdot
               end for
               Sample a batch of latent variables \{z^{(i)}\}_{i=1}^m \sim p(z).
11:
               \theta \leftarrow \operatorname{Adam}(\nabla_{\theta} \frac{1}{m} \sum_{i=1}^{m} -D_{w}(G_{\theta}(z)), \theta, \alpha, \beta_{1}, \beta_{2})
12:
13: end while
```

#### 13

<sup>13</sup>https://arxiv.org/pdf/1704.00028.pdf <sup>14</sup>https://arxiv.org/pdf/1701.07875.pdf

## Wasserstein GAN (Arjovsky et al. 17'<sup>15</sup>)



Figure 2: Different GAN architectures trained with different methods. We only succeeded in training every architecture with a shared set of hyperparameters using WGAN-GP.

<sup>15</sup>https://arxiv.org/pdf/1704.00028.pdf

## Wasserstein GAN (Arjovsky et al. 17'<sup>16</sup>)



Figure 3: CIFAR-10 Inception score over generator iterations (left) or wall-clock time (right) for four models: WGAN with weight clipping, WGAN-GP with RMSProp and Adam (to control for the optimizer), and DCGAN. WGAN-GP significantly outperforms weight clipping and performs comparably to DCGAN.

<sup>&</sup>lt;sup>16</sup>https://arxiv.org/pdf/1704.00028.pdf

#### Large scale WGAN (Karras et al. 18'<sup>17</sup>)



Figure 1: Our training starts with both the generator (G) and discriminator (D) having a low spatial resolution of  $4 \times 4$  pixels. As the training advances, we incrementally add layers to G and D, thus increasing the spatial resolution of the generated images. All existing layers remain trainable throughout the process. Here  $\boxed{N \times N}$  refers to convolutional layers operating on  $N \times N$  spatial resolution. This allows stable synthesis in high resolutions and also speeds up training considerably. One the right we show six example images generated using progressive growing at  $1024 \times 1024$ .

<sup>17</sup>https://arxiv.org/abs/1710.10196

# Large scale WGAN (Karras et al. $18'^{18}$ )



Figure 10: Top: Our CELEBA-HQ results. Next five rows: Nearest neighbors found from the training data, based on feature-space distance. We used activations from five VGG layers, as suggested by Chen & Koltun (2017). Only the crop highlighted in bottom right image was used for comparison in order to exclude image background and focus the search on matching facial features.

<sup>18</sup>https://arxiv.org/abs/1710.10196