









### Where is Deep Learning used?

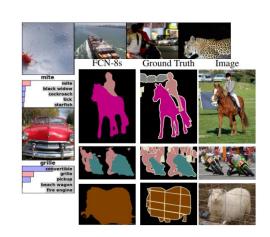
**Digit Recognition** 

Object Classification Segmentation

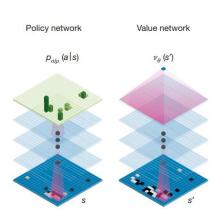
**Image Captioning** 

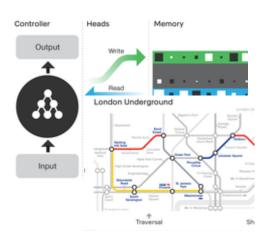
Gameplay Al Translation

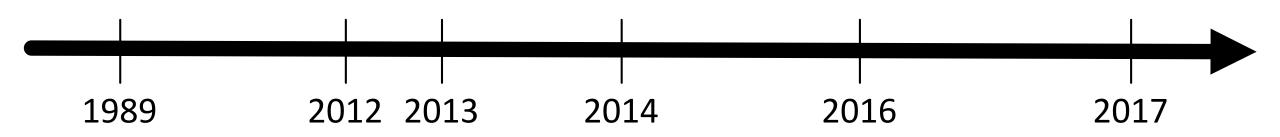
Neural Computers
Routing















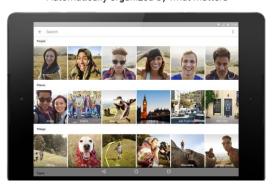


### Why Scale Up?

- Enormous amounts of data
  - MSCOCO: 19 GB
  - ImageNet (1k): 180 GB
  - ImageNet (22k): A few TB
  - Industry: Much larger
- Large neural network architectures
  - 100-200 layers deep today, ~100M-2B parameters
- Faster prototyping
  - Training time: 10s of hours to days (and weeks)



Automatically organized by what matters













## **Neural Networks**

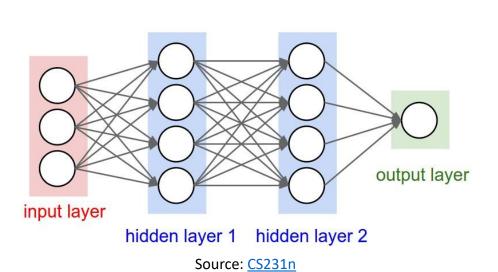


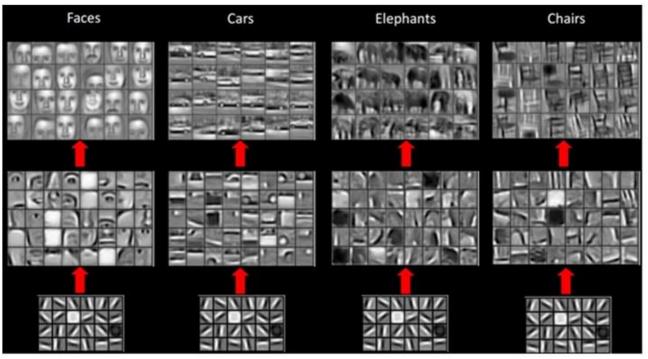




#### **Neural Networks**

- Modeled after the human brain
- CNNs repeatedly perform convolutions and nonlinearity operations





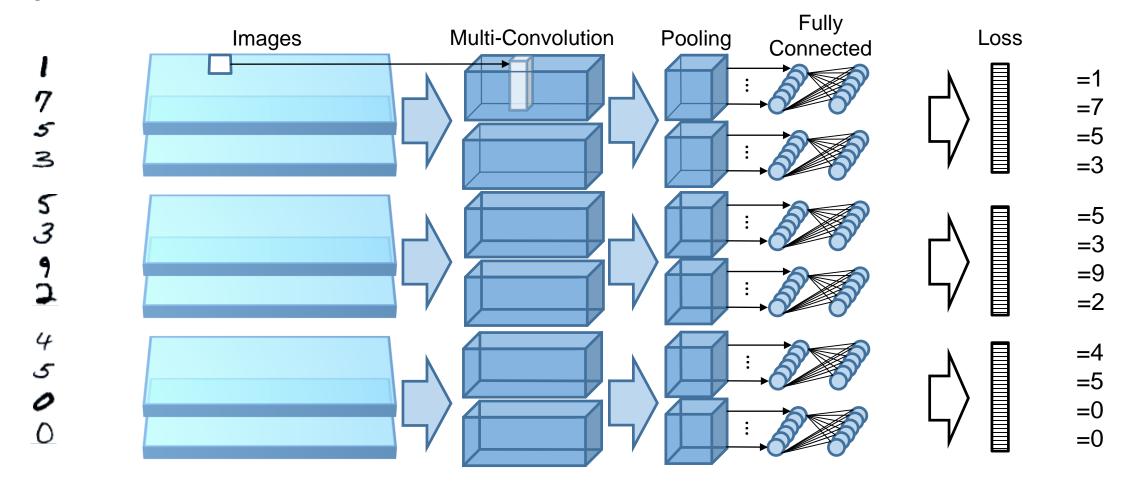
Source: Lee et al. <u>"Unsupervised Learning of Hierarchical Representations with Convolutional Deep Belief Networks"</u> (CACM 2011)







### **Simple CNN Architecture**



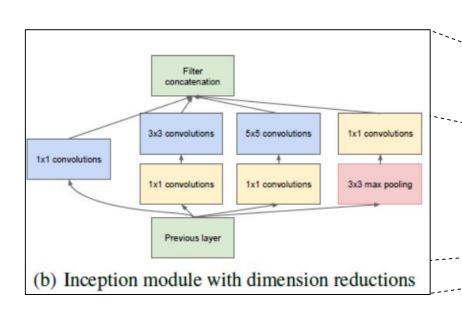


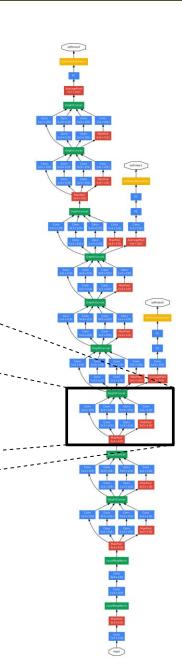




### GoogLeNet [Szegedy et al., 2014]

- ~6.8M parameters
- 22 layers deep





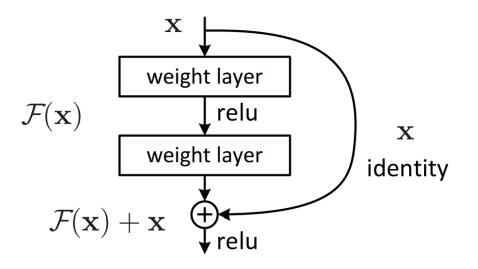


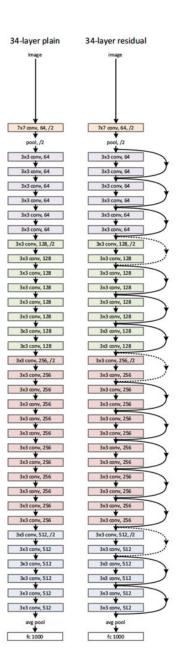




### ResNet [He et al., 2016]

- ~2.35M parameters
- 152 layers deep







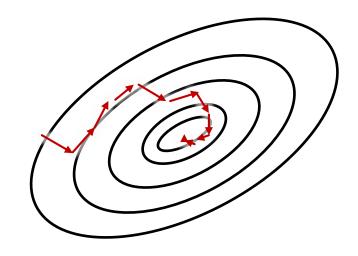




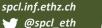
#### **Stochastic Gradient Descent**

- Gist: Improve network weights using samples from a labeled dataset
- Algorithm:
  - Initialize neural network weights  $(W_0)$
  - For t in iterations:
    - **Sample** *b* images from dataset (*B*)
    - Compute loss  $L_t(W_{t-1}, B)$
    - **Update** weights using gradients and update rule g:  $W_t = g(W_{t-1}, \nabla L_t(W_{t-1}, B), [hyperparameters ...])$
- $\nabla L_t(W, B)$  is an average direction of the gradient over a mini-batch of size b:

$$\nabla L_t(W, B) = \frac{1}{b} \sum_{i=1}^b \nabla \ell(W; (x_i, y_i))$$



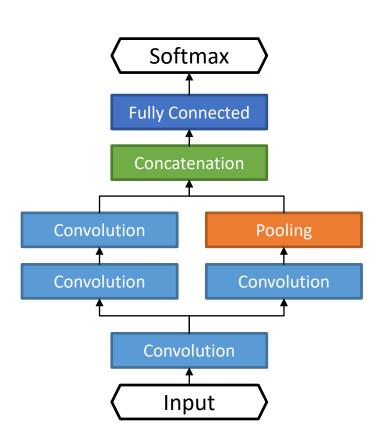






### **Backpropagation Algorithm**

- A CNN is a Directed Acyclic Graph (DAG)
- At each layer in backpropagation, derivatives are estimated w.r.t.:
  - Layer parameters (if necessary)
  - Data (chain rule)

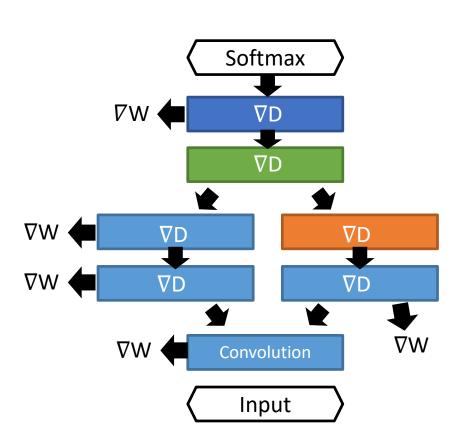






### **Backpropagation Algorithm**

- A CNN is a Directed Acyclic Graph (DAG)
- At each layer in backpropagation, derivatives are estimated w.r.t.:
  - Layer parameters (if necessary)
  - Data (chain rule)
- Additional memory storage required for training:
  - D+W+ **∇D+ ∇W**



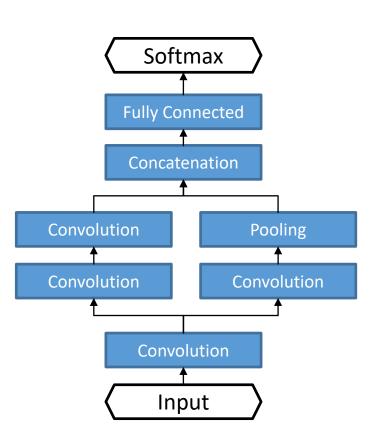






- Choice of Algorithm
- Parallelism

- Distributed Computing
- Hardware Architectures



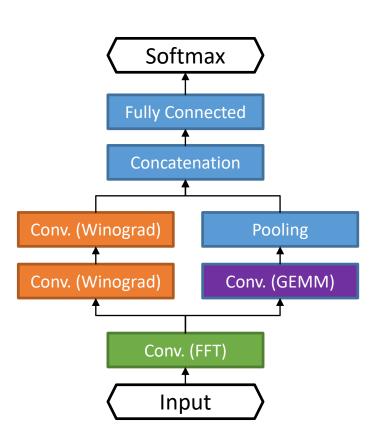






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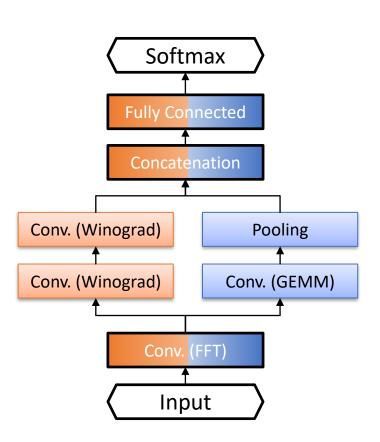
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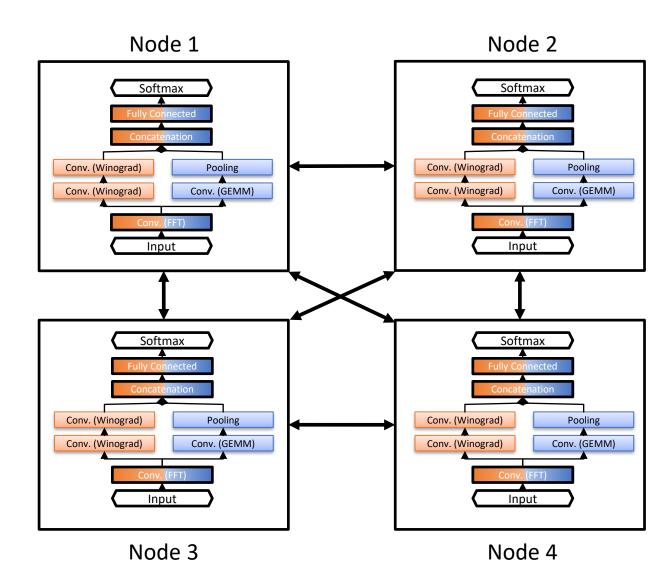




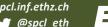


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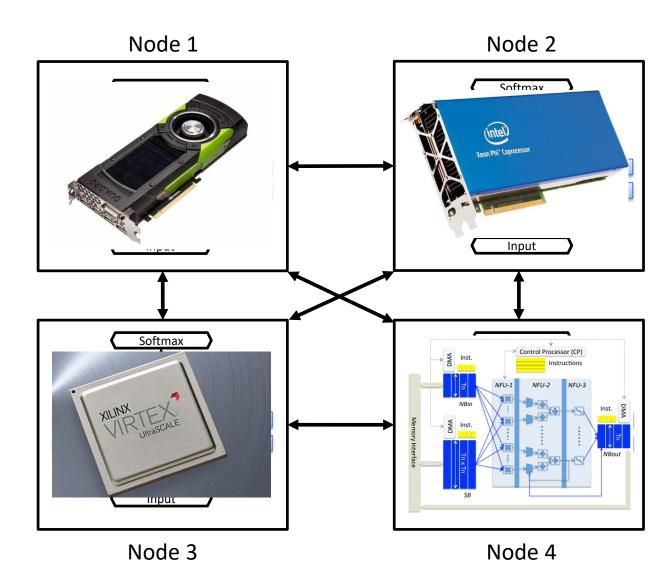






- Choice of Algorithm
- Parallelism

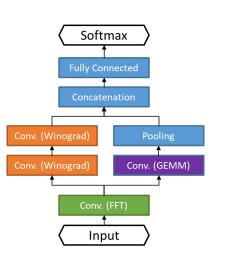
- Distributed Computing
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# **Algorithms**







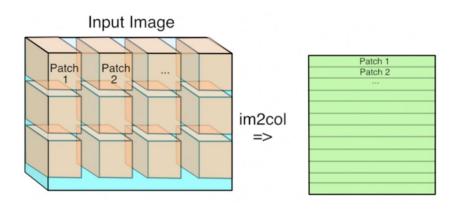


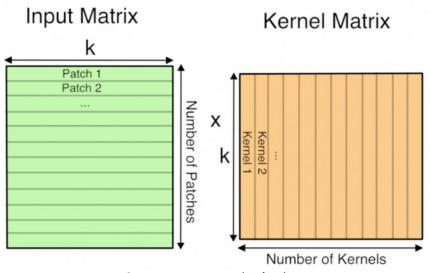
### **Convolution Algorithms**

Most computationally-intensive layer

$$out(x,y)^{f_o} = \sum_{f_i=0}^{N_{if}} \sum_{k_x=0}^{K_x} \sum_{k_y=0}^{K_y} w_{f_i,f_o}(k_x,k_y) *in(x+k_x,y+k_y)^{f_i}$$

- Can be performed directly, or:
  - Via matrix multiplication (im2col)
     [Chellapilla et al., 2006]
  - Via Winograd convolution [Lavin and Gray, 2016]
  - In Fourier domain





Source: Pete Warden's Blog

#### im2col convolution



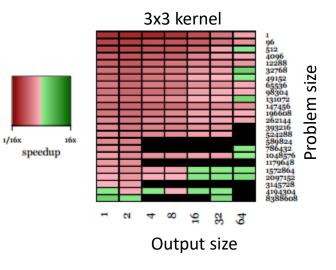


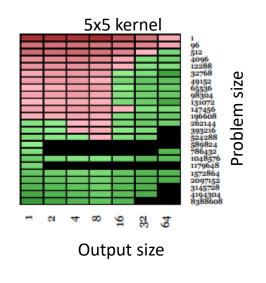
#### **Convolution in Fourier Domain**

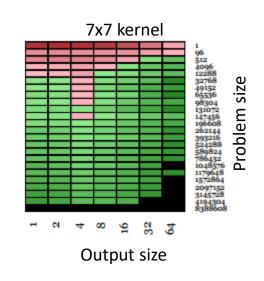
Convolution can be computed using FFT [Mathieu et al., 2014]:

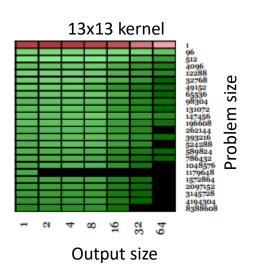
$$y_{(s,j)} = \sum_{i \in f} x_{(s,i)} \star w_{(j,i)} = \sum_{i \in f} \mathcal{F}^{-1} \left( \mathcal{F}(x_{(s,i)}) \circ \mathcal{F}(w_{(j,i)})^* \right)$$

The larger the convolution kernel, the better the performance [Vasilache et al., 2015]:













### **Sacrificing Accuracy for Performance**

- Half-precision (16-bit floating point) [Gupta et al., 2015]
  - Memory is stored in 16-bit format
  - Computations are performed in 32-bits
  - Uses Stochastic Rounding:

$$Round\left(x,\left\langle \mathtt{IL},\mathtt{FL}\right\rangle \right) = \begin{cases} \left\lfloor x\right\rfloor & \text{w.p. } 1 - \frac{x - \left\lfloor x\right\rfloor}{\epsilon} \\ \left\lfloor x\right\rfloor + \epsilon & \text{w.p. } \frac{x - \left\lfloor x\right\rfloor}{\epsilon} \end{cases} \end{cases}$$

*Goal:* Preserve  $\mathbb{E}(Round(x, \langle IL, FL \rangle)) = x$ 

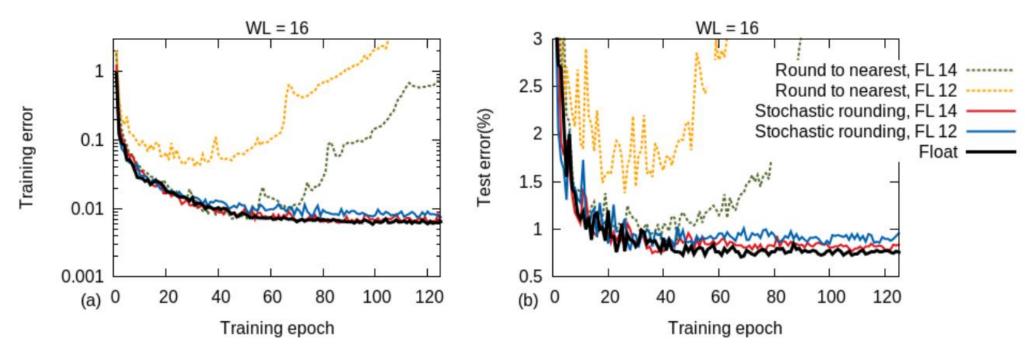






### **Sacrificing Accuracy for Performance**

Results on MNIST with LeNet:



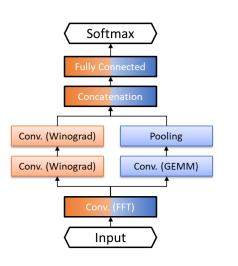
WL=Word Length (bits)
FL=Fractional Length (bits)







## **Parallelism**

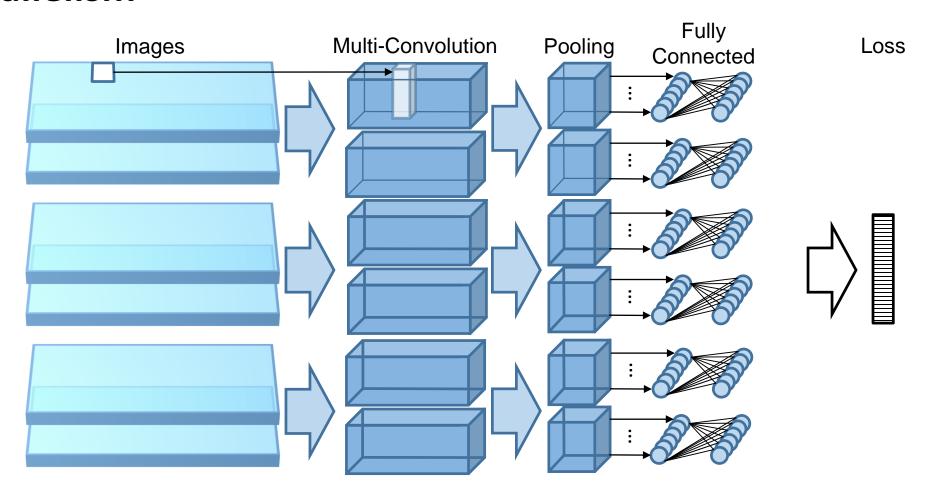








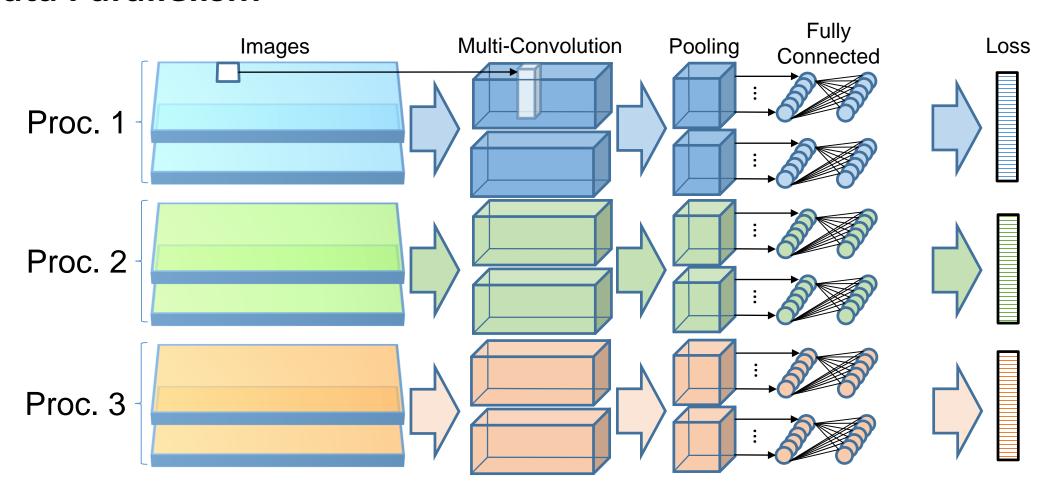
#### **Data Parallelism**







#### **Data Parallelism**









### **Data Parallelism**

- ✓ Good for forward pass (independent)
- ✓ Backpropagation requires all-to-all communication only when accumulating results

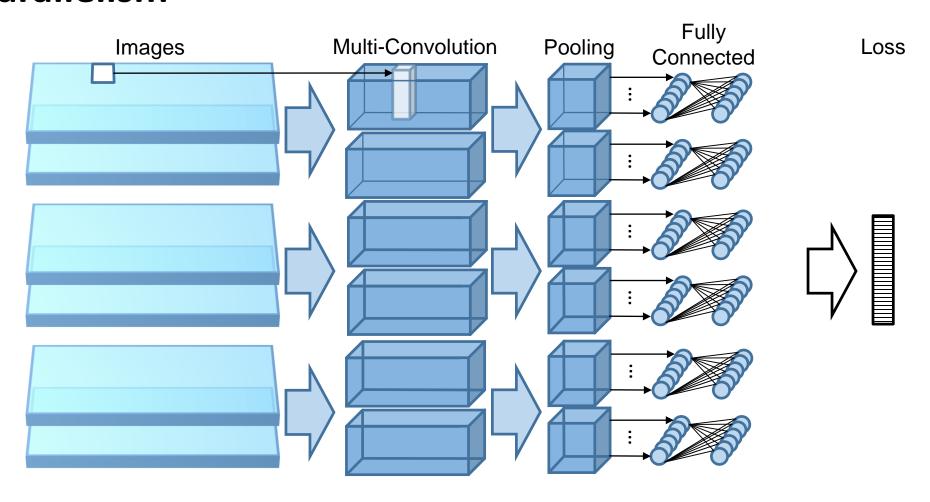
× Requires allocation of all parameters on each processor







#### **Model Parallelism**

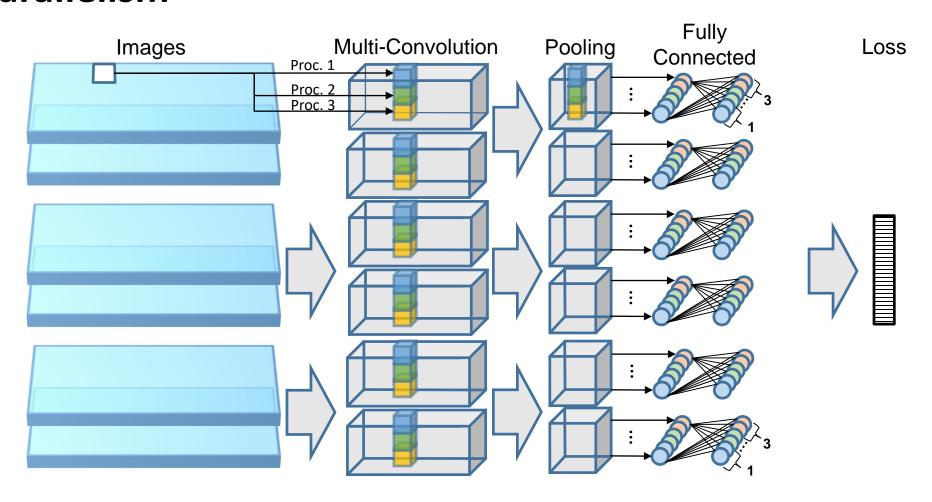








#### **Model Parallelism**









### **Model Parallelism**

- ✓ Parameters can be divided across processors
- × Mini-batch has to be copied to all processors
- × Backpropagation requires all-to-all communication every layer







### **Hybrid Data/Model Parallelism**

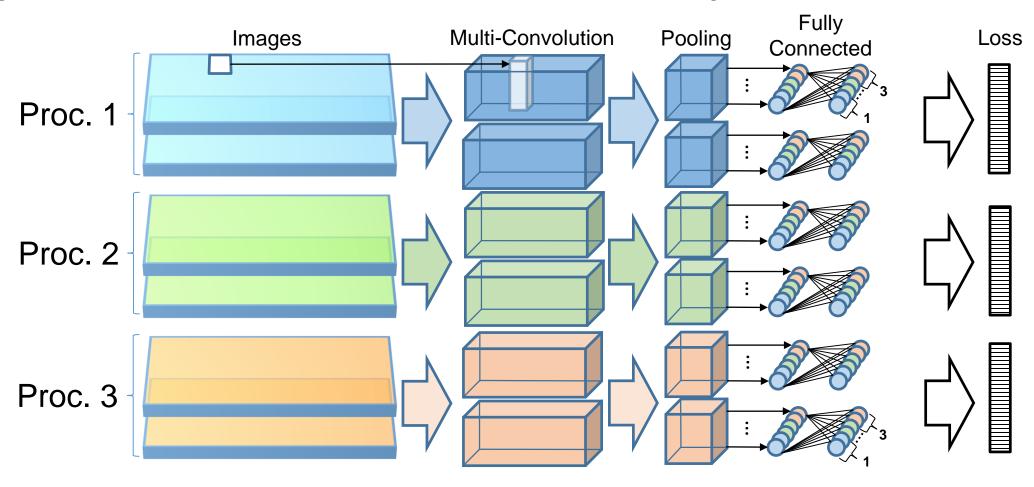
 Conjecture[Krizhevsky, 2014]: Most of the computations are performed in the convolutional portion, most of the parameters are stored in the fully connected portion

 Proposed Solution: Use data parallelism on convolutional portion and model parallelism on the FC portion





### Hybrid Data/Model Parallelism [Krizhevsky, 2014]









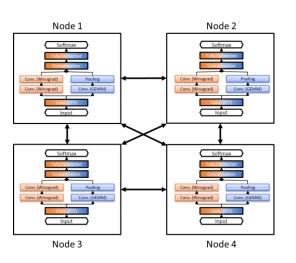
### **Hybrid Data/Model Parallelism Results**

AlexNet, ILSVRC 2012:

GPUs	Batch size	Top-1 error	Time	Speedup
1	(128, 128)	42.33%	98.05h	1x
2	(256, 256)	42.63%	50.24h	1.95x
2	(256, 128)	42.27%	50.90h	1.93x
4	(512, 512)	42.59%	26.20h	3.74x
4	(512, 128)	42.44%	26.78h	3.66x
8	(1024, 1024)	43.28%	15.68h	6.25x
8	(1024, 128)	42.86%	15.91h	6.16x



# **Distributed Computing**





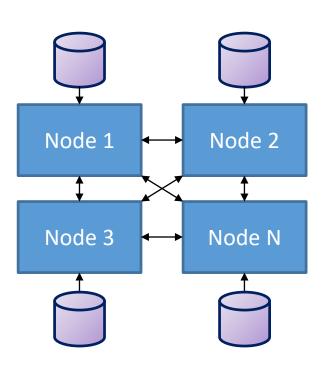




### **Distributed Deep Learning**

- Runs on a computer cluster
- Each node runs partially autonomously
- Inter-node communication from time to time

- Best result is gathered from the nodes
- Training data can be split to per-node "shards"



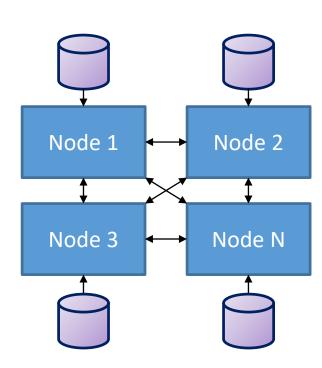






### **Distributed Deep Learning – Opportunities**

- Increased memory:
  - More data
  - More parameters
- Fault tolerance
  - Protection against node crashes
- Improved stochasticity



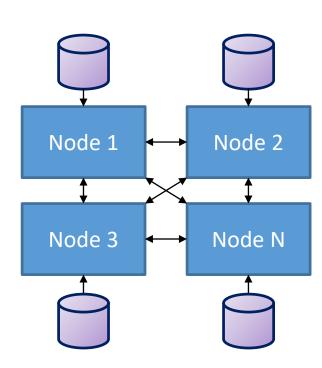






### **Distributed Deep Learning – Determining Factors**

- Computational independence
- Communication efficiency
- Network congestion
- Load balancing
- Points of failure



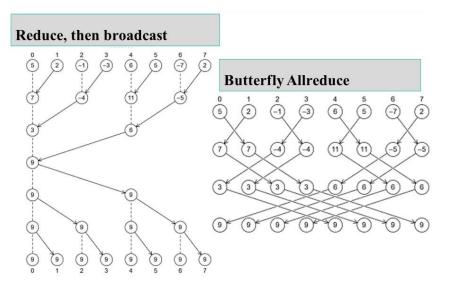






### **Distributed Synchronous SGD**

- Communication step is added to the algorithm:
  - Initialize neural network weights  $(W_0)$
  - For t in iterations:
    - Sample b images from dataset (B)
    - Compute loss  $L_t(W_{t-1}, B)$
    - Synchronize weights across workers
    - Update weights using gradients and update rule g



Source: Basics of Message-Passing

- The step requires all nodes to have the same data ( $\Sigma \nabla W$ )
  - This collective operation is also called AllReduce
- Different ways to implement, depending on message size and network topology



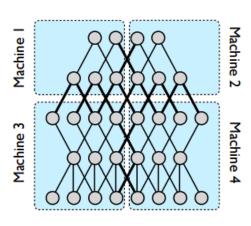


### **Distributed Deep Learning – DistBelief**

- Distributed learning infrastructure used at Google [Dean et al., 2012]
- Each model replica has the same parameters, but optimizes different data
  - Replicas are divided among several machines
- Two distributed optimization schemes for training:
  - Online Downpour SGD
  - Batch Sandblaster LBFGS



Handles slow and faulty replicas







### Asynchronous SGD – HOGWILD!

- To achieve coherency in distributed SGD, nodes must synchronize w.r.t. parameters:
  - Each thread draws a random example i from the training data.
    - Acquire a lock on the current state of parameters θ.
    - Thread reads  $\theta$ .
    - Thread updates  $\theta \leftarrow (\theta \alpha \nabla L(f_{\theta}(x_i), y_i))$ .
    - Release lock on  $\theta$ .
- *HOGWILD!* [Niu et al., 2011] removes this synchronization:

Each thread draws a random example i from the training data.

- Thread reads current state of  $\theta$ .
- Thread updates  $\theta \leftarrow (\theta \alpha \nabla L(f_{\theta}(x_i), y_i))$ .

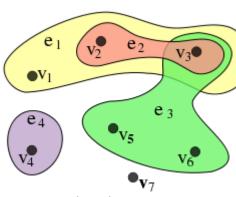






### Asynchronous SGD – HOGWILD!

- HOGWILD!:
  - Proven to converge in sparse problems
  - Provides near-linear scaling
  - Assumes shared-memory architecture (e.g., multicore CPUs)
- Formulates ML problems as hypergraphs G = (V, E) where:
  - $w^* = \operatorname{argmin}_w f(w) = \operatorname{argmin}_w \sum_{e \in E} f_e(w_e)$
  - Each hyperedge  $e \in E$  represents subsets of [n]
  - $w_e$  is reduced to coordinates in e



Source: Wikipedia

#### Algorithm 1 HOGWILD! update for individual processors

- loop
- 2: Sample e uniformly at random from E
- Read current state  $x_e$  and evaluate  $G_e(x_e)$
- 4: for  $v \in e$  do  $x_v \leftarrow x_v \gamma G_{ev}(x_e)$
- 5: end loop

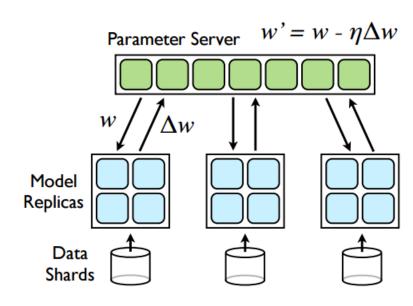






### **Distributed Deep Learning – Downpour SGD**

- Relaxation of HOGWILD! for distributed systems
- Algorithm:
  - Divide training data into subsets and run a replica on each subset
  - Every  $n_{fetch}$  iterations, fetch up-to-date parameters from server
  - Every  $n_{push}$  iterations, push local gradients to server
- Note that parameter shards may be "out-of-sync"



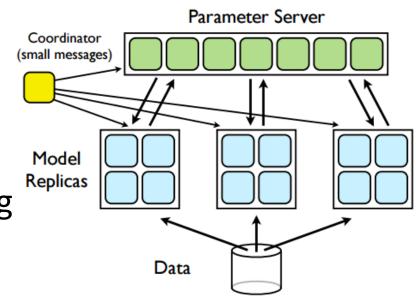






### **Distributed Deep Learning – Sandblaster LBFGS**

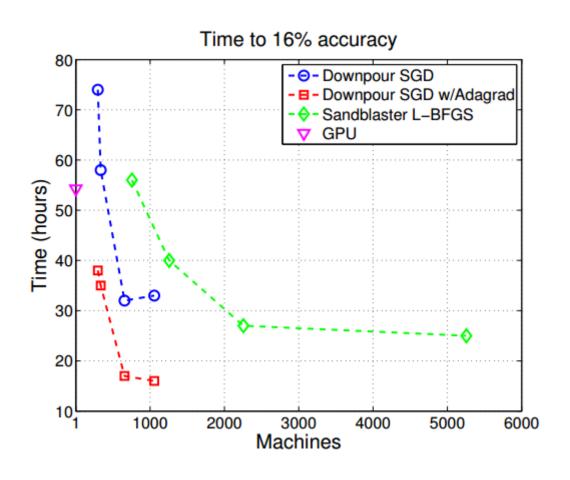
- Coordinator process issues commands (dot product, scaling, multiplication, etc.) to slave nodes, each processing a different parameter shard
- Communication is sparser
  - Most of the information is stored locally
  - Coordinator messages are small
  - Slaves fetch parameters at the beginning of each batch, send gradients once in a while for fault tolerance
- Employs computation replication and load balancing
  - Nodes that finish their job get more jobs
  - If one node is slow, additional nodes get the same job







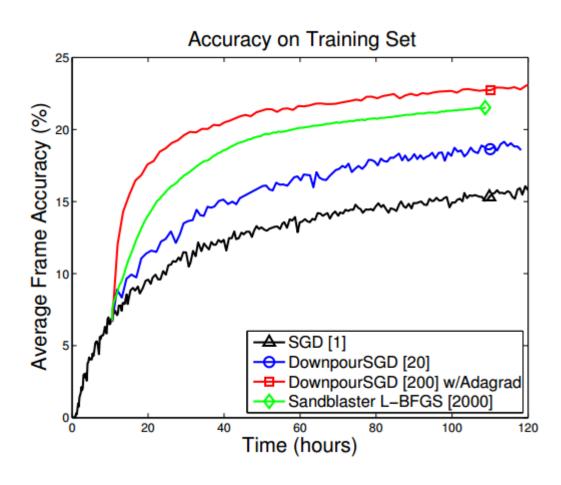
### **DistBelief Results – Time**

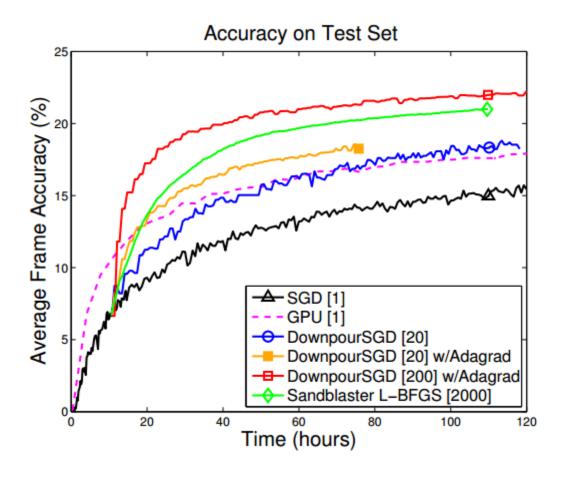






### **DistBelief Results – Accuracy**





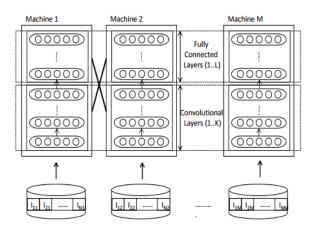


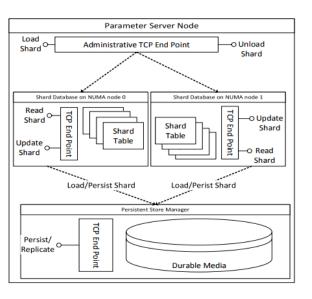




### **Project Adam**

- Extends DistBelief with system-level support
  - Fast data serving mechanisms (e.g., with augmentation)
  - Better heterogeneous system management
  - Parameter server node optimization
- Bottom-up communication redesign
  - Control message, data message separation
  - Inter-node communication reduction
  - Weight differences are sent instead of weights
- Only system to train ImageNet22k

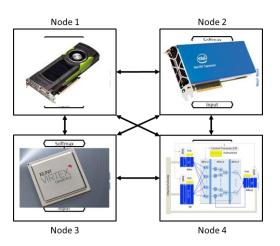








# Hardware







### **Specialized Hardware**

- GPU
  - Thousands of cores, massively parallel (5-14 TFLOP/s per card)
  - Multi-GPU nodes further increase training performance (using data/model parallelism)
  - Drawback: Hard to program efficiently. Solution: Specialized libraries (CUDNN)
- FPGA
  - Specialized for certain operations (e.g. convolutions)
  - Drawbacks: Even harder to program
- Convolutional Processing Units

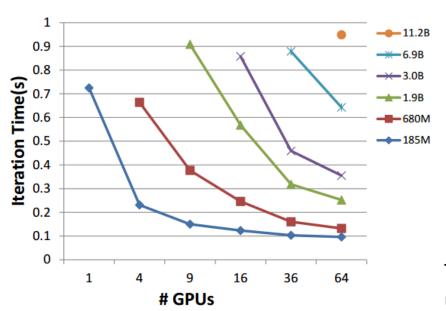






### **Deep Learning with GPUs**

- A distributed GPU-based system[Coates et al., 2013] was shown to run DistBelief-scale problems (1000 machines) with 3 multi-GPU nodes
- 3 tiers of concurrency: GPU, model parallelism, nodes

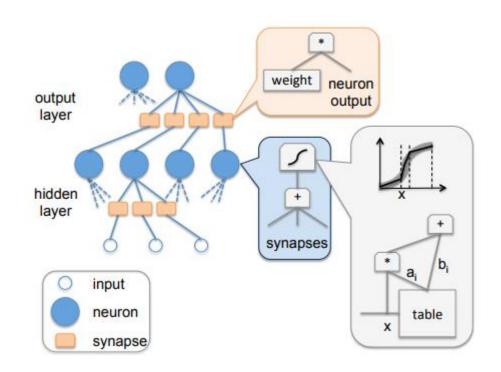


Time for a single mini-batch gradient update of a sparse autoencoder

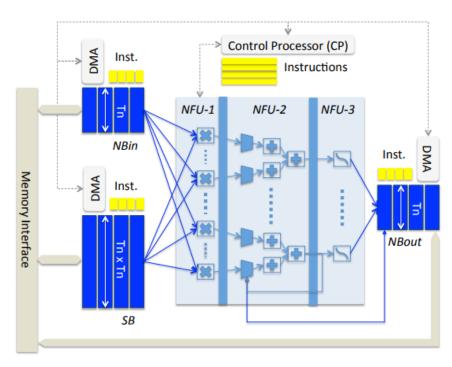


## **Specialized Hardware**

Two approaches:



Mapping neurons to hardware



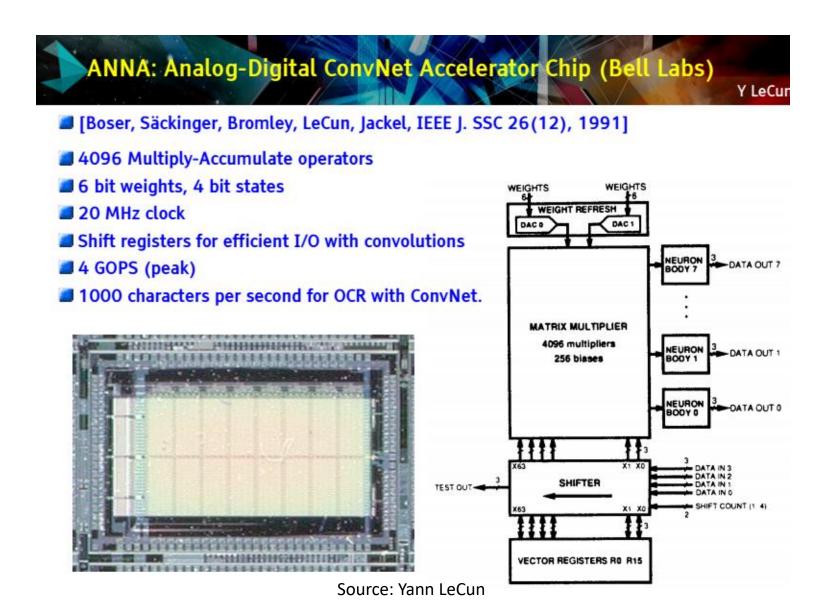
**Custom processing elements** 







### **Specialized Hardware – ANNA**



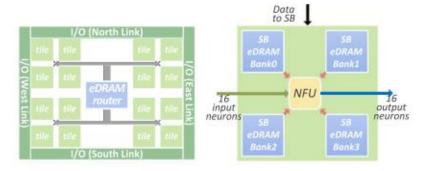




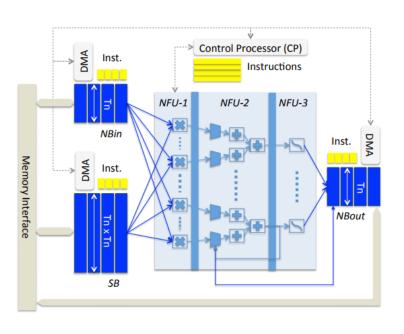


### **Specialized Hardware – DianNao**

 Instead of building the neural network on a circuit, creates a Neural Function Unit (NFU)



- Operations reflect the different layers
- Each operation (conv., pooling, FC) takes up to three stages at computations (or less)
  - Computation
  - Reduction
  - Activation









### **Conclusions**

- Many acceleration opportunities
- Architectures keep changing, and with them new techniques arise
- Algorithms can be modified (to some extent)
  - Proven "shortcuts" can be taken









### **Questions?**