Large Scale Machine Learning

Distributed learning - Part 2

Previous session: Synchronous Distributed SGD

- Trade-off between
 - Compute time
 - Communication time
 - Synchronisation time
- Reducing communication cost
 - Exploiting sparsity
 - Compression
 - But Increasing computation
- Works reasonably well in practice
 - Can also give a good initial solution to be fine tuned with more complex methods

In Spark

```
// Load training data in LIBSVM format.
val data = MLUtils.loadLibSVMFile(sc, "data/mllib/sample_libsvm_data.txt")
// Split data into training (60%) and test (40%).
val splits = data.randomSplit(Array(0.6, 0.4), seed = 11L)
val training = splits(0).cache()
val test = splits(1)
// Run training algorithm to build the model
val model = new LogisticRegressionWithSGD()
                  .setNumClasses(10)
                  .run(training)
// Compute raw scores on the test set.
val predictionAndLabels = test.map { case LabeledPoint(label, features) =>
    val prediction = model.predict(features)
    (prediction, label)
```

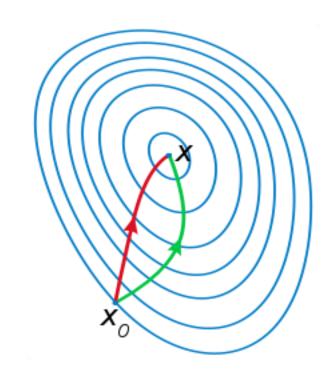
This session

- Decreasing required number of iterations
 - Second order methods, Newton / LBFGS
 - ADAM
- Decreasing communication cost
 - Tree reduce, ring reduce
- Decreasing synchronization cost
 - Asynchronous updates!
- Beyond data parallelism

Second order optimization methods

 We want to speed up our convergence in another way

 Using the "curvature" information can help a lot



Newton's method (1 dimension)

• In one dimension:

Quadratic approximation of function L to minimize:

$$L(x) = L(x_t) + (x - x_t) L'(x_t) + (x - x_t)^2 L''(x_t)/2 + o((x - x_t)^2)$$

Step minimizing local approximation: $x_{t+1} := x_t - (L''(x_t))^{-1} \cdot L'(x_t)$

Very good convergence properties

Newton's method

Taylor expansion of L:

$$L(x) = L(x_t) + (x - x_t) \cdot \nabla L(x_t) + (x - x_t) \cdot \nabla^2 L \cdot (x - x_t) + o((x - x_t)^2)$$

Step minimizing local approximation:

$$X_{t+1} := X_t - (\nabla^2 L(X_t))^{-1} \cdot \nabla L(X_t)$$

Hessian matrix

- Coefficients $\frac{\partial^2}{\partial x_i \partial x_j} L(x)$
- Size n²
- Very good convergence properties on convex problems
- Hessian is of size n²

Quasi-newton methods

- Instead of: $dx = -(\nabla^2 L(x_t))^{-1} \cdot \nabla L(x_t)$
- Approximate $(\nabla^2 L(x))^{-1}$ by some matrix B_t
- $dx = -\alpha B_t \nabla L(x_0)$

Select step size α with Line search

- Several methods, usually estimating \boldsymbol{B}_t from previous gradients and steps
- L-BFGS

Quasi-newton methods: generic

- Until convergence
 - Compute update direction $(B_n \approx H_n^{-1})$ $\Delta \theta = -B_n g_n$
 - Line search for learning rate

$$\alpha \leftarrow \min_{\alpha \ge 0} L(\theta_n - \alpha \Delta \theta)$$

Update parameters

$$\theta_{n+1} \leftarrow \theta_n - \alpha \Delta \theta$$

Store the parameters and gradient deltas

$$g_{n+1} \leftarrow \nabla L(\theta_{n+1})$$

$$\Delta \theta_{n+1} \leftarrow \theta_{n+1} - \theta_n$$

$$\Delta g_{n+1} \leftarrow g_{n+1} - g_n$$

Update inverse hessian approximation

$$B_{n+1} \leftarrow QuasiUpdate(B_n, \Delta\theta_{n+1}, \Delta g_{n+1})$$

BFGS

Update B with a rank two matrix, of the form

$$B_{n+1} \leftarrow B_n + auu^T + bvv^T$$

- Limited memory BFGS or L-BFGS for short
 - Perform the update without actually materializing B matrix and performing an explicit matrix vector multiplication
 - Can be achieved by storing the latest few values and gradients

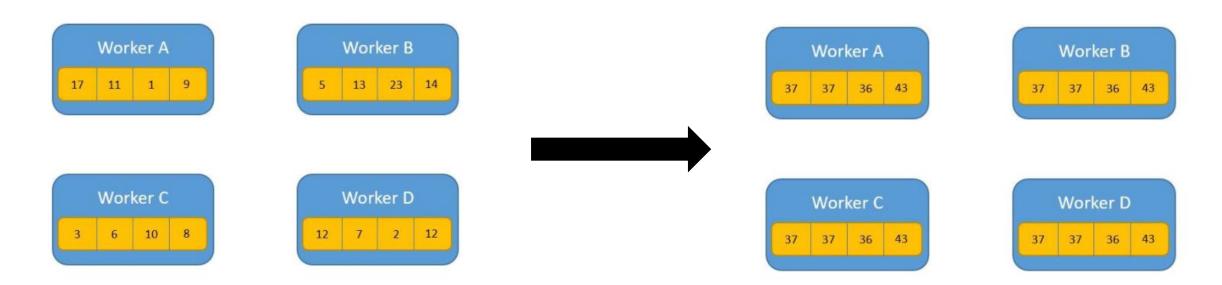
ADAM

```
alpha = 0.01 # Stepsize
           beta1 = 0.9
 3
           beta2 = 0.999
 4
           epsilon = 1e-8
 5
           epsilon = epsilon
 6
           m = v = t = 0
 8
       def AdamStep(gradient):
 9
           m = beta1 * m + (1.0 - beta1) * gradient
           v = beta2 * v + (1.0 - beta2) * gradient * gradient
10
11
           mhat = m / (1.0 - beta1**(t+1))
12
           vhat = v / (1.0 - beta2**(t+1))
13
           step = alpha * mhat / (np.sqrt(vhat) + 1e-8)
14
           step = step / ( 1+ decay * t )
15
           t += 1
16
           return step
17
18
     # ... on each example ...
19
     w = w - AdamStep(gradient)
```

AllReduce

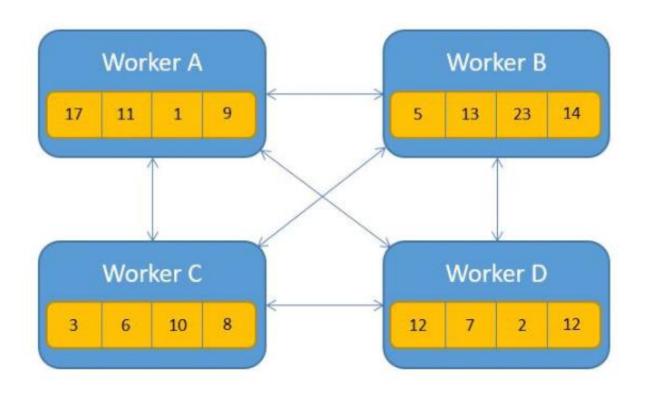
Local gradients on the workers We need to:

- Sum all local gradients
- Send the sum to all workers

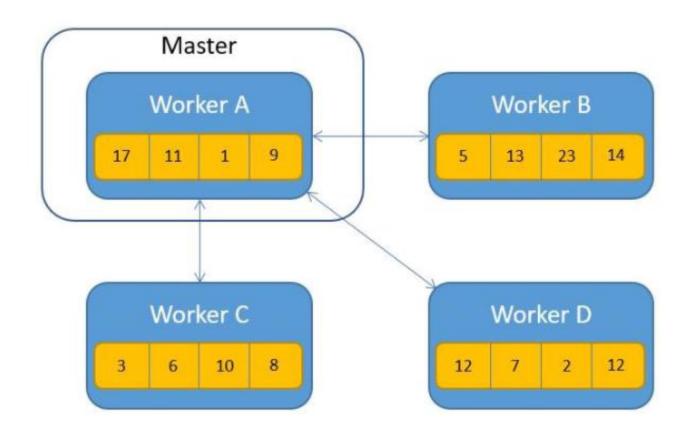


https://towardsdatascience.com/visual-intuition-on-ring-allreduce-for-distributed-deep-learning-d1f34b4911da

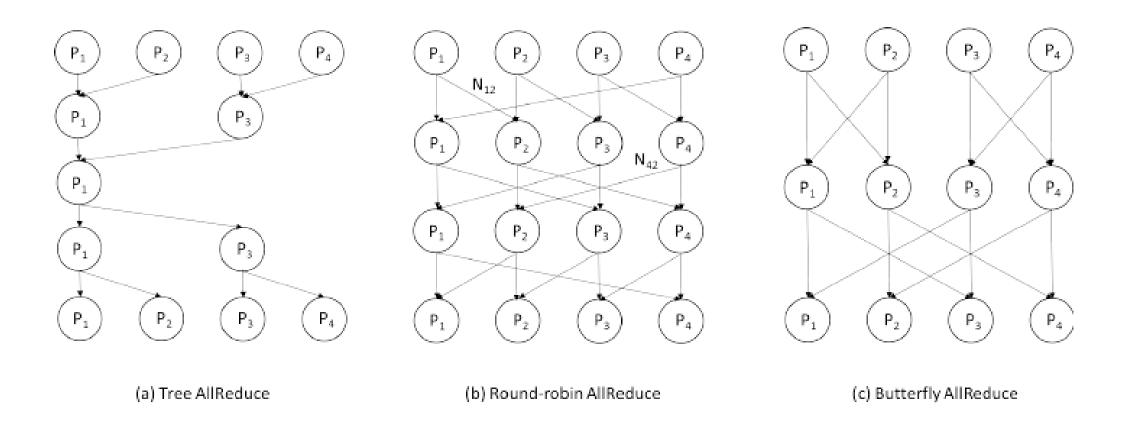
AllReduce: all to all communication

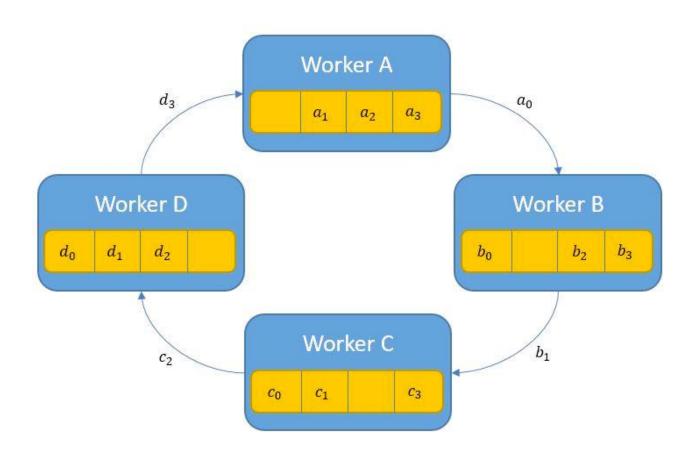


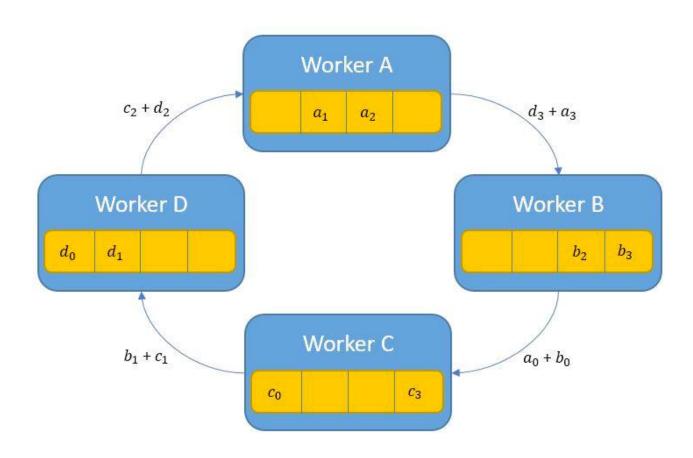
AllReduce: Reduce then Broadcast

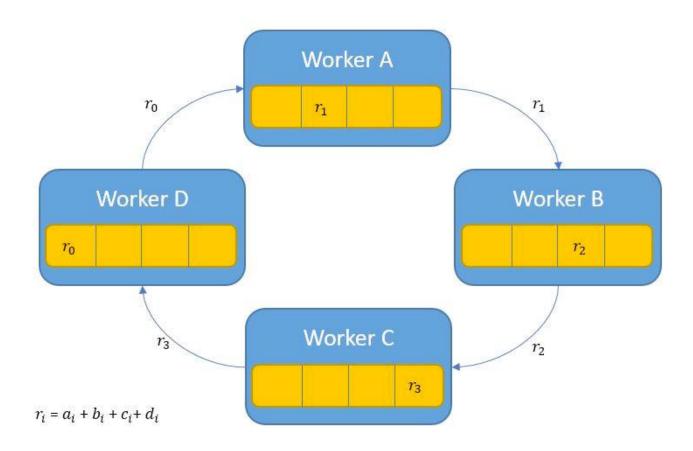


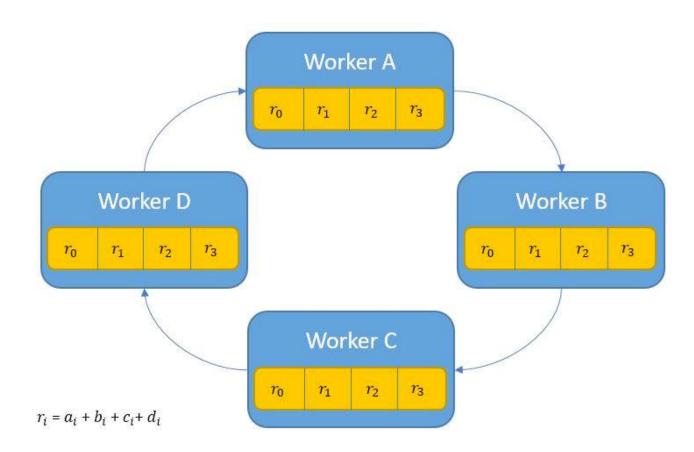
AllReduce



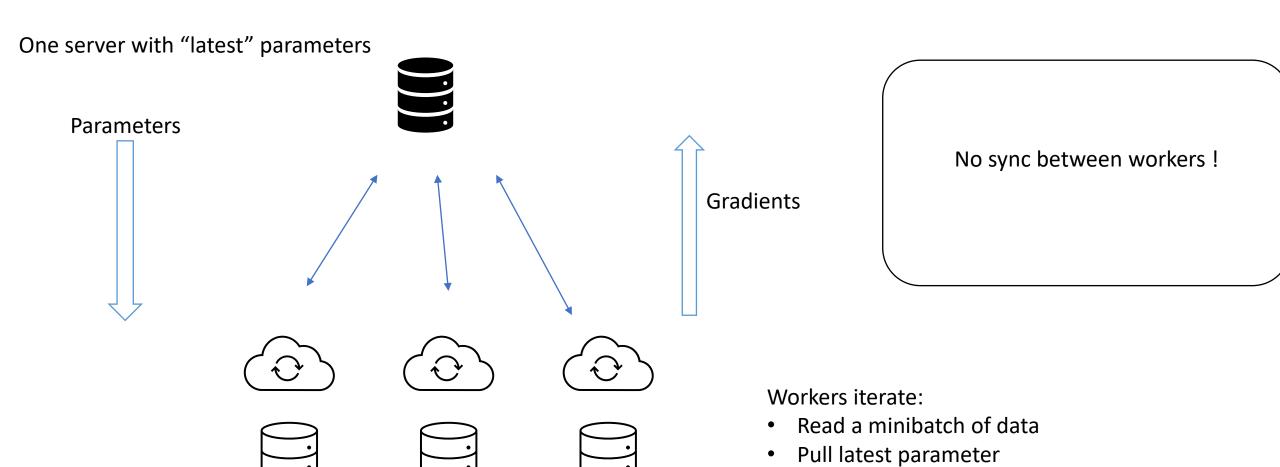








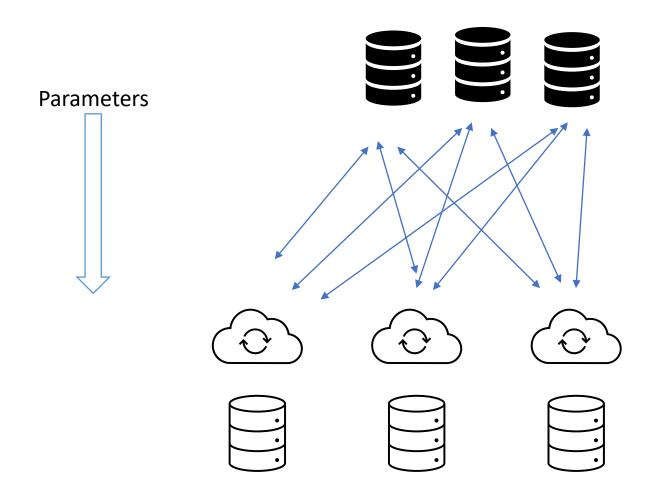
Parameter server to avoid synchronization



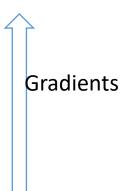
Compute the gradient

Send parameters updates

Bottleneck: Bandwidth on the master?

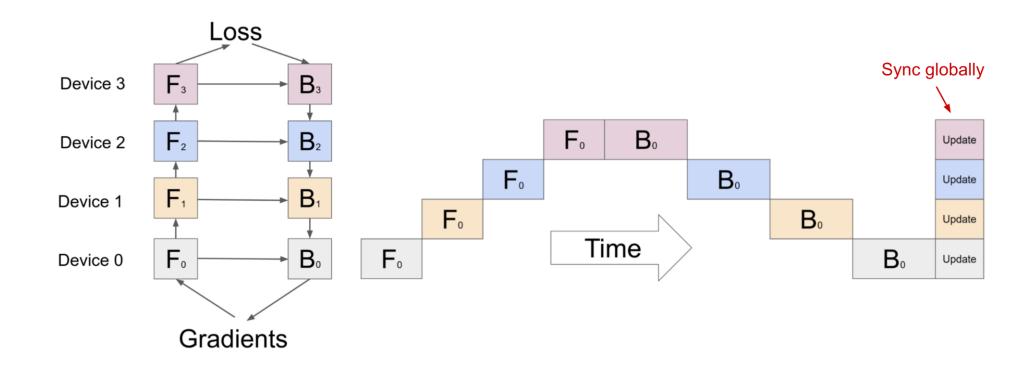


Split the parameters on several servers!

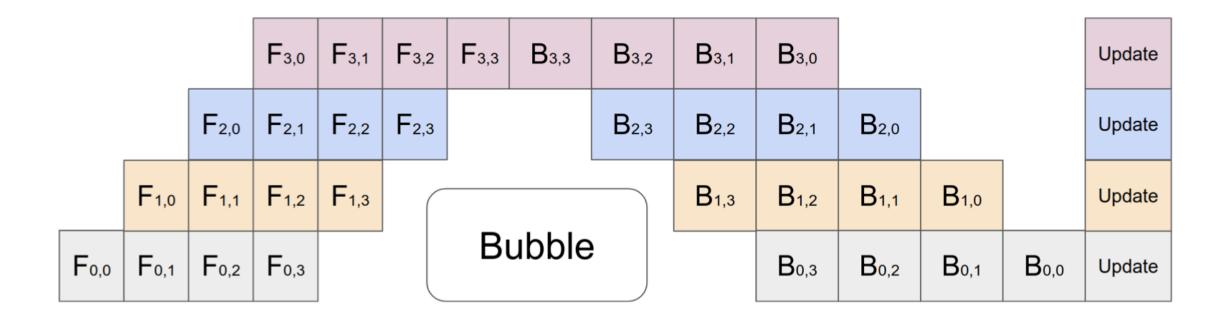


Beyond data parallelism

- Very large models: too big to fit on a single worker
- => Requires to split a model among several workers, eg one layer / worker



Pipeline parallelism model parallelism + data parallelism



Asynchonous pipeline parallelism

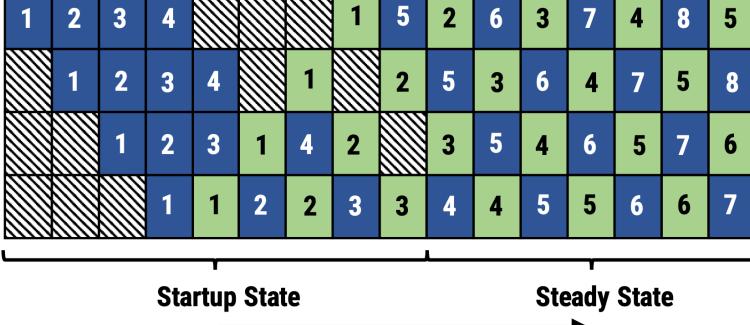
Pipedream https://cs.stanford.edu/~matei/papers/2019/sosp_pipedream.pdf

Machine 1

Machine 2

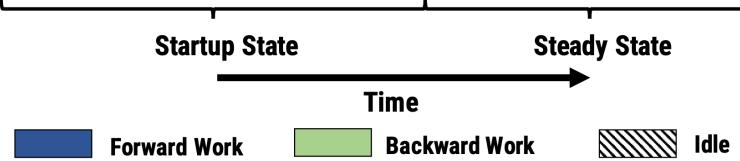
Machine 3

Machine 4



Complication:

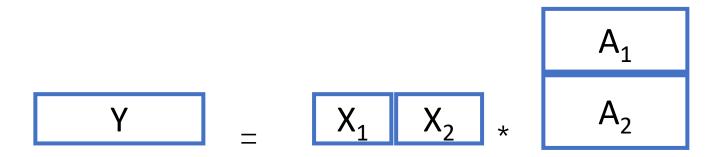
Ensuring the same version of the weights are used in forward and backward passes



Tensor parrallelism

• When a single layer is too large (eg large input embedding layer)

Typically, decomposing a matrix multiplication by block



- compute each block on different workers
- requires a synchronization to gather the blocks

Conclusion

- Distributed Machine Learning is about trade-offs
 - Communication VS computation cost
- For simple models (like Logistic Regression), a synchronous approach works well
 - Exploit sparsity
 - Use more complex optimization schemes
- There are several ways to distribute and aggregate computation
 - Centralized synchronous or asynchronous model, AllReduce ...