TTIC 31230 Fundamentals of Deep Learning, winter 2019 Backpropagation Problems

Problem 1: Backprogation through a ReLU linear threshold unit. Consider the computation

$$y = \sigma(w^{\top}x)$$
$$\ell = \mathcal{L}(y)$$

for $w, x \in R^d$ with $\sigma(z) = \max(z, 0)$ (the ReLU activation) and for $\mathcal{L}(y)$ an arbitrary function (a loss function). Let w_i denote the *i*th component of the weight vector w. Give an expression for $\frac{\partial \ell}{\partial w_i}$ as a function of $\frac{d\mathcal{L}(y)}{dy}$.

Solution: There are various correct ways of writing the answer. The following corresponds to a backpropagation computation.

$$\frac{d\ell}{dy} = \frac{d\mathcal{L}(y)}{dy}$$

$$\frac{d\ell}{dw_i} = \frac{d\ell}{dy} \frac{dy}{dw_i} = \frac{d\ell}{dy} x_i \mathbf{1} \left[w^\top x \ge 0 \right]$$

Problem 2: Backpropagation through softmax. Consider the following softmax.

$$Z[b] = \sum_{j} \exp(s[b, j])$$

$$p[b, j] = \exp(s[b, j])/Z[b]$$

An alternative way to compute this is to initialize the tensors Z and p to zero and then execute the following loops.

for
$$b, j$$
 $Z[b] += \exp(s[b, j])$
for b, j $p[b, j] += \exp(s[b, j])/Z[b]$

Each individual += operation inside the loops can be treated independently in backpropagation.

(a) Give a back-propagation loop over += updates based on the second loop for adding to s.grad using p.grad (and using the forward-computed tensors Z and s).

Solution: For b, j s.grad $[b, j] += p.\text{grad}[b, j] \exp(s[b, j])/Z[b]$

(b) Give a back-propagation loop over += updates based on the second equation for adding to Z.grad using p.grad (and using the forward-computed tensors s and Z).

Solution: For b, j $Z.\operatorname{grad}[b] = p.\operatorname{grad}[b, j] \exp(s[b, j])/Z[b]^2$

(c) Give a back-propagation loop over += updates based on the first equation for adding to s.grad using Z.grad (and using the forward-computed tensor s).

Solution: For b, j s.grad[b, j] += Z.grad $[b] \exp(s[b, j])$

Problem 3: Optimizing Backpropagation through softmax. Show that the addition to s.grad shown in problem 2 can be computed using the following more efficient updates.

for
$$b, j$$
 $e[b] = p[b, j]p.\operatorname{grad}[b, j]$
for b, j $s.\operatorname{grad}[b, j] += p[b, j](p.\operatorname{grad}[b, j] + e[b])$

Solution: The updates for problem 1 can be written as

for
$$b$$
 $Z.\operatorname{grad}[b] = \sum_{j} -p.\operatorname{grad}[b, j] \exp(s[b, j])/Z[b]^2$

$$= \left(\sum_{j} -p[b, j]p.\operatorname{grad}[b, j]\right)/Z[b]$$

$$= e[b]/Z[b]$$

$$\begin{array}{lll} \text{for } b,j & s. \text{grad}[b,j] & = & p. \text{grad}[b,j] \exp(s[b,j])/Z[b] + Z. \text{grad}[b] \exp(s[b,j]) \\ & = & p. \text{grad}[b,j] \left(\exp(s[b,j])/Z[b]\right) + e[b] \left(\exp(s[b,j])/Z[b]\right) \\ & = & p[b,j](p. \text{grad}[b,j] + e[b]) \end{array}$$

This formula shows how hand-written back-propagation methods for "layers" such as softmax can be more efficient than compiler-generated back-propagation code. While optimizing compilers can of course be written, one must keep in mind the trade-off between the abstraction level of the programming language and the efficiency of the generated code.

Problem 4. Backpropogation through batch normalization. Consider the following set of += statements defining batch normalization where all computed tensors are initialized to zero.

For
$$b, j$$
 $\mu[j] += \frac{1}{B} x[b, j]$
For b, j $s[j] += \frac{1}{B-1} (x[b, j] - \mu[j])^2$
For b, j $x'[b, j] += \frac{x[b, j] - \mu[j]}{\sqrt{s[j]}}$

Give backpropagation += (or -=) loops for computing $x.\operatorname{grad}[b,j]$, $\mu.\operatorname{grad}[j]$, and $s.\operatorname{grad}[j]$ from $x'.\operatorname{grad}[b,j]$. The loops should be given in the order they are to be executed.

Solution:

For
$$b, j$$
 $x.grad[b, j]$ += $\frac{x'.grad[b, j]}{\sqrt{s[j]}}$
For b, j $\mu.grad[j]$ -= $\frac{x'.grad[b, j]}{\sqrt{s[j]}}$
For b, j $s.grad[j]$ -= $\frac{1}{2}(x[b, j] - \mu[j])s[j]^{-3/2} x'.grad[b, j]$
For b, j $x.grad[b, j]$ += $\frac{2}{B-1} (x[b, j] - \mu[j])s.grad[j]$
For b, j $\mu.grad[j]$ -= $\frac{2}{B-1} (x[b, j] - \mu[j])s.grad[j]$
For b, j $x.grad[b, j]$ += $\frac{1}{B} \mu.grad[j]$

Problem 5. Writing framework code. Consider a function $c: R^d \times R^s \to R^s$, in other words a function that takes a vector of dimension d and a vector of dimension s and yields a vector of dimension s. Given a sequence of vectors x_0, x_2, \ldots, x_T with $x_t \in R^d$ we can define a sequence of vectors h_0, h_1, \ldots, h_T by the equations

$$h_0 = c(x_0, 0)$$

 $h_t = c(x_t, h_{t-1}) \text{ for } 1 \le t \le T$

When the function c is defined by a neural network the resulting network mapping x_1, \ldots, x_T to h_0, \ldots, h_T is called a recurrent neural network (RNN). a. In the educational framework EDF we work with objects where each object has a value attribute and a gradient attribute each of which have tensor values where the value tensor and the gradient tensor are the same shape. Each object is assigned a value in a forward pass and assigned a gradient in a backward pass. Suppose that we are given an EDF procedure CELL which takes as arguments a parameter object Phi and two EDF objects X and H where the value attribute of the object X is a d-dimensional vector and the value attribute of the object H is an s-dimensional vector. A call to the procedure CELL(Phi, X, H) returns an EDF object whose value attribute is computed in a forward pass in some possibly complex way from the value attributes of Phi, X and H. Given a sequence X[] of EDF objects whose value attributes are d-dimensional vectors, and an EDF object ZERO representing the constant s-dimensional zero vector, write a procedure for constructing the sequence of EDF objects representing h_1, h_2, \ldots h_T as defined by the above RNN equations. Your solution can be in Python or informal high level pseudo code.

Solution: We can use the equations given as the definition of the computation graph if we replace c in the equations with the function CELL. In the folloing code CELL is a class parameter packages and the call CELL() creaates a fresh parameter package on each call. A recuarive solution can also be given.

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\begin{split} &X = list() \\ &H = list() \\ &H[0] = CELL(Phi(), X[0], ZERO) \\ &for \ t \ in \ range(1,T) \\ &H[t] = CELL(Phi, X[t], H[t-1]) \end{split}
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b. Deep learning systems generally make extensive use of parallel computation for training. How does the parallel running time of an RNN computation graph scale with the length T?

Solution: The parallel running time is proportional to T. RNNS are fundamentally serial and this is a problem. RNNs have recently been largely replaced by the transformer architecture.