1 Introduction

Imagine you are my professor. Maybe you actually were my professor, in which case you may already be sweating before I say any more. The subject matter is Neural Networks. You draw an illustration on the board with a node’s inputs, and its output via a transfer function.

“Now this transfer function can be almost anything. Typically it would be something like the hyperbolic tangent, which looks like this.

But it has to be a non-linear function. If it’s linear, i.e. of the form \( y = mx + b \), then observe that the entire layer is a linear function. And so the entire network is just a linear function of linear functions; itself a linear function. We could just compute an equivalent single-layer network, and we know that it could only fit linear functions, which is insufficient for most problems.”

Then I raise my hand. The speed with which I raise it, and the subtle forward pose of my arm suggests that I want to pluck an abstract idea from the whiteboard and pervert it. You know this look, and you’re reluctant to call on me. But no other students are asking questions. You must call on me.

“Tom.” It’s more like a statement then a question. It includes the tone of spoken punctuation that, if it could, ends the entire conversation before it begins.

“OK but, when we implement this on a computer we’ll use some approximation of numbers, like floating point. So the specific sequence of additions and multiplications will matter. It’s not actually equivalent to rearrange them to a single layer because you don’t have distributivity, commutativity, etc.”

That was about 20 years ago. The world will not let us stop thinking about neural networks. And so this question has been on my mind for a long time. Just to be clear, the professor was right: This is not an important question. Theoretically I am right, but for practical purposes it probably does not matter. But I like to work at the intersection of Theory and Impractice. We can make it matter by doing a lot of work. And then I will continue to be right theoretically, but also more right because it will only matter for most practical purposes.

So this paper is an exhaustive exploration of what we can do with just floating point addition and multiplication by constants (scaling). You should only be able to make lines, but I’ll demonstrate that due to rounding error, you can absolutely use “linear” transfer functions in neural networks. Machine learning is not the only field with a proclamation that some function must be “non-linear,” so we’ll look at a few of those as well. There will of course be several hearty digressions. By studying these functions we’ll see that they are almost arbitrarily rich, and conclude with a demonstration of their completeness in the field of Plumbing.

2 A refresher on neural networks

Let’s repeat the professor’s lesson. This section is easily skippable if you are a plucky student who thinks they already know everything. At a high level, a Neural Network is a way of implementing a numeric function (takes a bunch
of numbers as input, and gives a bunch of numbers as output. The network consists of a number of layers, where the first layer is the input and the last layer is the output. Each layer is an array of nodes. Here is a simple three-layer network with some of the nodes labeled:

![Image of a neural network diagram]

The numbers that fill in each layer are its activations (here some of these values are labeled a, b, ...). Each layer’s activations are computed from (just) the previous layer. Looking at the bold portion in the example, the value of e is given as

\[ e = TF(w_0a + w_1b + w_2c + w_3d + \text{bias}) \]

The multiplicative weight \((w_i)\) and additive bias \((\text{bias})\) parameters are learned during the training of the neural network, but just become constants when using the neural network to compute its output.

\(TF\) is the transfer function, which is of particular interest in this project. Classically, the transfer function was some kind of sigmoid. The \(\tanh\) function pictured in the introduction is a good example of a sigmoid. The intuition behind this is that, thinking about a node as some kind of neuron, the neuron “fires” (activates) with some probability. This probability gets higher as its input values get larger, but can’t be higher than 1. Note that weights can be negative, so upstream neurons can have an inhibitory effect. In fact it is frequently useful for neurons to “negatively fire” (outputting \(-1\)). The \(\tanh\) function clamps the result symmetrically to \((-1, 1)\) rather than a probability.

**Differentiability.** Another important property of the transfer function is that it be differentiable, because the stochastic gradient descent algorithm used to train neural networks needs to be able to move along some error-reducing gradient, and back-propagate errors to earlier layers. This gradient is just the derivative of the function.

**What transfer functions ought to exist?** We used to think that these saturating transfer functions were ideal. But this turns out to be wrong, especially for internal (“hidden”) layers. Transfer functions don’t need to produce probabilities, and they can have unbounded range.

A wide variety of functions will work, including extremely simple ones. The most popular transfer function in 2023 is the “rectified linear unit,” which looks like this:

\[ x \leq 0 ? 0 : x \]

This one is extremely easy to implement \((x < 0 ? 0 : x)\), is fast and seems to work very well, possibly because its derivative is significant (one) on the entire positive side. (In contrast, sigmoids tend to get “stuck” because of their saturating behavior; their derivatives become nearly zero when activations are high.) Note that it is not actually differentiable (discontinuity at zero) but “for all practical purposes” it is differentiable.

The (only?) apparently essential quality of the transfer function is that it be non-linear. If it is instead of the form \(TF(x) = mx + b\), then any activation \(a\) is also just a linear function of the previous layer, as linear functions of linear functions (weighted sum) are linear. This causes the entire network to be a linear function. It is well known that a linear function “cannot” represent some other simple functions, such as XOR.

\[ \forall m, n, b. \ XOR(x, y) \approx mx + ny + b \]

This means that with a linear transfer function, a neural network could never learn even a simple function like XOR. Many problems we want to learn are in fact much more complicated.

### 3 A fine terminological issue

My smart math friend Jason refers to a function like \(f(x) = mx + b\) pejoratively as “high school linear.” Depending on what class you’re in, this may formally be an affine function because of the bias term \(b\). Here I use “linear” to mean a polynomial of degree \(\leq 1\). If you wanna perjorate me as being in high school, so be it.

**The Rules.** To be precise, we will allow addition and scaling by constants. When we have a “linear” function of multiple variables, these variables can be individually scaled and added, but not multiplied. So for a function like \(f(x, y, z) = x + 3y - 2z + 4\) is allowed, as is anything mathematically equivalent to it (like \(f(x, y, z) = 2x + 4 + 2y - 2z - x + y - 0\). \(f(x, y, z) = xy + z\) is not permitted.

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1 In these contexts, a linear function must obey \(f(0 \times x) = 0 \times f(x)\), so it must be zero at zero.
4  Half-precision IEEE-754 floating point

In this project we’ll abuse floating point inaccuracy to create “linear” functions (only using floating point addition and scaling) that are not lines. For this reason, we prefer to have a numerical system that is less accurate. In floating point, inaccuracy comes from the fact that not all numbers are representable (due to finite precision) and the result of an operation is always rounded to a representable number. IEEE-754 floating point [1] comes in different “spice levels,” with “32-bits” being “float” and “64-bits” being “double.” Although spice levels as low as 3 bits make sense [27], 8-bit (“mild”) is quite common in machine learning. Usually the reason to prefer half precision is that it uses less memory, and so your GPU can store networks that are twice as big in RAM. For this project we will also use half precision, and we will be happy to save RAM, but more happy that its precision is low and so it is practical (although silly) to achieve significant rounding error. Another important reason to choose half precision is to make the pun in the title.

A half precision float is 16 bits: One sign bit, five bits for the exponent, and 10 bits for the mantissa. Like all IEEE-754 formats, there is much more precision (more values are representable) near zero (Figure 1). Once you get to 1024, only integers are representable. From 2048 to 4096, only even numbers are representable. 65504 is the largest finite number, and up here, only multiples of 32 are available.

Some CPUs have native support for half-precision IEEE-754, but typically via non-standard intrinsics or compiler flags. Since people using half-precision are usually doing so in the interests of performance, many configurations will “help” you by performing practical but incorrect optimizations. This is similar to what happens when enabling --ffast-math, which stands for Final Fantasy AST Math, meaning that the abstract syntax tree of your program will be manipulated using fantasies about Math that do not apply to IEEE-754, and your Final result can be arbitrarily different. For the ideas in this paper to work, --ffast-math is prohibited. And it will be slow!

Rather than deal with non-standard stuff, I found a nice library called half.h [29] that implements IEEE-754 compliant half-precision in portable C++. I use this throughout the project and it matches the behavior of my GPU. I recommend it for similar hijinks.

_origins of imprecision. Floating point does have many perversions, but many programmers come to believe all sorts of dangerous superstitions about it. One idea is that floating point is somehow always inexact, and so that you always have to check that two numbers are equal “within some epsilon” [23]. This may work “in practice” but it is actually pretty sloppy. Floating point imprecision is not random, nor is it constrained to a fixed epsilon. Operations are defined much more usefully: Each one computes the mathematically correct value, and then rounds (according to the “rounding mode”) to the nearest representable value. That’s it. One consequence of this is that you can get the exact result of 32-bit multiplication by doing 64-bit multiplication and then rounding to 32 bits. This also means that the rounding error from a single operation can be as large as the gap between representable numbers: Up to 32 for half-precision. But it also means that operations whose results can be exactly represented have no error; for example adding integral half values less than 512 will always give an exact integer result, which can be compared using ==. We will use this later in Section 7.1. It is neither necessary nor sufficient compare for “equality” with some “epsilon.”

Rounding. IEEE-754 supports multiple rounding modes, like “round-to-zero,” and “round-to-infinity” (always round in the positive direction). Throughout this paper we use “round-to-nearest,” which is also the typical default (e.g. for C++11 expressions evaluated at compile time, it always uses round-to-nearest)\(^2\). Similar results are likely attainable for the other rounding modes, as well as hypothetical rounding modes such as “round away from nearest,” but I have not explored this.

Getting some nonlinearity. All transfer functions implemented with floating point have a finite range. For our experiments with neural networks, we will focus on transfer functions that map values in \([-1, 1]\) to values in \([-1, 1]\). Almost half (48.4\%) of floating point values are in this interval and this is a typical nominal range for activations in neural networks.

\(^2\)There is seldom reason to change the rounding mode, and since it is a stateful act, you’re asking for it if you do. But the round-to-negative-infinity and round-to-positive-infinity modes are useful for interval arithmetic, which is arguably the only truly reasonable way to use floating point. What you do is represent numbers as intervals (low and high endpoints) that contain the true value, and then perform each calculation on both endpoints. For computations on the low endpoint, you round down, and symmetrically for the high endpoint. This way, the true value is always within the interval, and you also know how much inaccuracy you have accumulated!
We only have two operations: Addition and scaling. Let’s see what kind of rounding error each of these gives us. First, addition. In order to get a function that takes values in $[-1, 1]$ to values in $[-1, 1]$, we want to first add a constant (giving us perhaps a large value) and then add a negative constant, bringing us back in range. For example, the constant 128 gives us the function

$$f(x) = x + 128.0 - 128.0$$

This is of course mathematically the same as $f(x) = x$ (the identity), but with half precision we get a function that looks like this.

Between 128 and 256, only multiples of 0.125 are representable. So for arguments in 0 to 1, the sum is rounded to one of the values 128.0, 128.125, 128.25, . . . 129. From 64 to 128, multiples of 0.0625 ($\frac{1}{16}$th) are representable. So from –1 to 0, we get 127.0, 127.0625, 127.125, . . . 128. Subtracting 128, all of the values are exactly representable, giving us $-1, -0.9375, . . . , -0.0625, 0, 0.125, . . . , 0.875, 1$.

The result is a step function, but whose resolution is twice as high for the negative range as the positive; had we added –128 and then added 128, we would have seen the opposite bias in resolution. We can easily see that this function is (computationally) non-linear despite being (mathematically) “linear.” This function is unlikely to be a good transfer function, because for one thing it does not have a good derivative: It’s zero most places (flat segments) except at the discontinuities, where it is undefined. We do test this approach (with the constant 64.0) later, though.

Scaling gives similar results. Consider

$$f(x) = x \times 100.0 \times (1.0/100.0)$$

In this project we never actually divide (although this would not violate linearity) since most floating point numbers have approximate multiplicative inverses, and many are exact. We just compute the reciprocal $\frac{1}{100} \cong 0.01000213623$ ahead of time and multiply by that constant. Here’s what that function looks like:

At this scale it appears linear, but it does have small imperfections (see zoomed region). The function is symmetric about zero, since multiplication will do the same thing to a positive number as it does to its negative counterpart. Here, the roundoff error differs with the magnitude. At inputs close to 1.0, the results of the first multiplication must round to the nearest multiple of 0.0625 (as in the additive example) but this error is scaled down by a factor of 100 when we multiply back to the $[-1, 1]$ range. So it is almost invisible. For inputs close to 0.0, the error approaches zero. The effect is complex and depends on the constant we multiply by. For example, if we multiply by a power of two, this only affects the exponent, and so the result is exact.

Is that it? Of course not! We can apply these operations in combination, and many times, to create more interesting functions. The best approach I found in this simple family is to repeatedly multiply the input by a number very close to one. Here’s what happens if you multiply the input by 0.99951171875 (which is the next number smaller than one, equal to $1 - \frac{1}{2048}$) five hundred times, and then scale back at the end:

$$f(x) = x \times (1 - \frac{1}{2048}) \times (1 - \frac{1}{2048}) \times \ldots 500 \text{ times} \times \frac{1}{13232421875}$$

I call this the grad1 function.

Multiplying 1.0 by ($1 - \frac{1}{2048}$) five hundred times in half precision yields 0.755859375 (mathematically it would be $1 - \frac{1}{2048})^{500} = 0.78333$, so there is significant accumulated error. We set $f(1.0) = 1.0$ by multiplying by the inverse of this constant, which is 1.3232421875.

Why does this result in the zig-zags? Multiplication by ($1 - \frac{1}{2048}$) affects numbers differently. For constants less than 6.1094760895 × 10⁻², the value is unchanged; we round back up to the original value. For all other finite inputs it produces a smaller value, but with rounding error that depends on the value. This error accumulates and becomes significant with many iterations (Figure 2). Unlike the previous functions, the output here is much smoother (it looks
Figure 2: How repeatedly multiplying by \( 1 - \frac{1}{2048} \) affects values in \([0, 1]\). The width of the image is the interval \([0, 1]\), with zero at the left.

**Top:** In the topmost row, we assign each pixel a hue so that we can track where those values go. For each pixel, we successively multiply by the constant and plot its color in its new \( x \) position, the move to the next row down. Note that the rainbow shrinks exponentially as expected, but not smoothly. The black line is 500 iterations.

**Bottom:** The accumulated error when iteratively multiplying by the constant and plot its color in its new \( x \) position, the move to the next row down. Note that the rainbow shrinks exponentially as expected, but not smoothly. The black line is 500 iterations.

piecewise-linear); in each of these segments its derivative is nondegenerate. Of course, this function is mathematically linear. It is equivalent to \( f(x) = x \times 1.036535 \).

So now we have a “good” candidate function, which we’ll call \( \text{grad}1 \). It is “good” in the sense that it is computation-ally non-linear despite being mathematically linear, so it may prove my professor wrong. On the other hand, it requires 501 floating point multiplications to compute, which is kind of slow. The “good” news is that since there are only 65536 16-bit values, we can easily just precompute any function for all possible half inputs, and store it in a table of 131072 bytes. This allows us to execute the function efficiently when performance is important, such as during training. (Table lookup is certainly not a mathematically linear operation, so when we require the computation to be linear for ideological purposes, we can perform the 501 multiplications and get the same result.)

**Differentiating.** Speaking of training, in order to train a neural network using stochastic gradient descent, we need to be able to evaluate the derivative of the transfer function at any point. We use that derivative to decide what direction to move the parameters (it gives us the “gradient” that we “descend”) as we propagate errors back through the network. There is an annoyance here, or if you like, an opportunity for a trick. We typically store the activation of each node, which is the output of the transfer function, but the derivative of a function is normally described in terms of the input (for example we say if \( f(x) = x^2 \) then \( f'(x) = 2x \)). We could store both the input and output for this step, or store only the input and recreate the outputs by running the transfer function. But the trick: We can compute the derivative as function of the output. For \( f(x) = x^2 \) we could say \( f'(f(x)) = 2\sqrt{x} \). Oops! That doesn’t actually work for \( x^2 \) because the square root could either be negative or positive, and the derivative is different depending on which one it is. In order for this trick to work, the transfer function has to be injective. Fortunately this is the case for the classic transfer functions, and this trick is well known so you don’t even need to do any math; you just look the function up.

For new transfer functions like \( \text{grad}1 \), we need to figure something out. This function does appear injective if we squint at it, although it is not really injective if you zoom way in: There are some distinct inputs that result in the same output due to rounding. But this is true for almost all floating-point functions already. I’ll be damned if I can come up with an analytic derivative for this thing, though. At best it would be some piecewise linear thing, requiring some table. Since our domain is only 16-bit, it is completely practical to just table the entire derivative (keyed by the output value, as we need). I do this programmatically. We do not want the derivative to reflect the step function that we see at very fine scales (the derivative should never be 0 for this function, for example), so I use a lowpass filter. The

\[ f(x_1) = f(x_2), f'(x_1) = f'(x_2). \] For the rectified linear unit, for example, all negative inputs are mapped to zero. But the derivative is also just zero in this entire region.
Since we need the derivative in terms of \( \text{grad1}'s \) output, the derivative is oriented along the \( y \) axis; each blue dot's \( x \) coordinate gives the derivative at the point on the black line that shares a \( y \) coordinate. It's an oscilloscope!

result looks good, oscillating between two different slopes as expected (Figure 3). The derivative is loaded into GPU memory during training and the table lookups are plenty fast.

4.1 Bonus digression: Downshift

Having freed myself from needing to “do math” in order to differentiate exotic functions, I pondered other weird transfer functions. For example, the rectified linear transfer function is very simple and works well, but is it the fastest possible transfer function that might work? It does involve a conditional, which naively implies comparison and branching (although probably most processors can do this with a conditional move). Because the floating point format is packed with fields that represent different things, many simple operations on its bits have interesting non-linear behavior. The most promising I found was a right shift by two places. It looks like this:

Shifting is about the cheapest possible thing a processor can do. Its behavior on floating point numbers is interesting:

Note the different regions for sign, exponent, and mantissa. The sign bit is shifted into the exponent, which means that the output is always non-negative (like the rectified linear function) and is non-linear (discontinuity at zero, as negative numbers have a much larger exponent than positive ones). Further nonlinearity comes from the exponential representation (shifts divide the exponent by four) and reinterpretation of exponent bits as mantissa bits. There is additional weirdness in the details. Shifting by two places is better than one, as it cannot produce Inf or NaN. We will also evaluate this transfer function, called downshift2, below.

Back to the main topic. I implemented all this as a modification of my custom neural network training and inference system, “Tom7Flow.” Tom7Flow is generally much worse than mainstream packages; it is based on deprecated OpenCL technology, is prone to divergence or stagnation during training due to naive choices of hyperparameters, etc. But it is at least well suited to silly experiments that take the form, “What if deep learning but worse?” such as the current exercise. In order to realize the idea completely, I modified the inference code to calculate with half-precision arithmetic (not just the transfer function). This means that the trained networks can be executed using only half-precision operations (and just addition and multiplication by constants). Unfortunately, while my GPU supports half-precision math natively, and OpenCL supports half-precision operations as an extension [??], this extension is somehow not supported (??) by my drivers, perhaps because OpenCL is so thoroughly deprecated. It does support half precision as a storage format, which allows you to write a full-precision float to a 16-bit value (rounding to half) or read a 16-bit half into a float (all half values can be represented exactly in full precision). So with this one operation it is straightforward to implement half-precision addition and scaling. You maintain the invariant that any float value is always exactly a half, and after you perform addition or multiplication, you round to half (by storing in a 16-bit memory location and reading it back). This definitionally produces the same results as the native operation.

I initially tried a version of training that worked entirely using half precision (network parameters are half, backpropagated errors and update values are half, etc.). This worked badly. It is ideologically unnecessary, as we just care about producing a final model that, during inference, only executes linear half-precision operations (but abuses floating point roundoff to do something interesting.) This network can be trained using non-linear techniques (and must anyway, since for example its computed derivative is not linear). So during training, calculations are done using full-precision floats, except for the forward step (where we round to half after every operation). In addition to being simpler, representing intermediate learned weights as floats seems to help training approach the final half values smoothly, avoiding stalls due to underflow.

\[4\] I also verified consistent results using the \texttt{half.h} software implementation. Many of the evaluation results quoted in the paper are actually executed on the CPU using this library.
4.2 Neural network experimental results

In order to evaluate this transfer function, I ran a suite of benchmark problems. For each problem, I compare the same network architecture (i.e. the number of layers, their connectivity, random initialization, etc.) but using different transfer functions.

The transfer functions are:

- **grad1**: The “linear” transfer function grad1 described above.
- **tanh**: The hyperbolic tangent function, which is a classic saturating (output is always in \((-1, 1)\)) sigmoid.
- **logistic**: The function \(1/(1+e^{-x})\), another classic sigmoid (but whose output is in \((0, 1)\)). Each operation is performed with half precision.
- **leaky relu**: The rectified linear unit, but with a small slope below zero: \(x < 0.0 \circ 0.1 \cdot x : x\). This is the function I usually prefer in practice; its advantage over the standard relu is that it does not “die” (zero propagated error) when its input is negative.
- **downshift2**: Interpreting the half-precision input as a 16-bit word, right shift by 2 places, then reinterpret as half.
- **plus64**: \(f(x) = x + 64 - 64\). This about the simplest function that has obvious rounding error. It only outputs 25 distinct values in \([-1, 1]\) so its derivative is degenerate; I use its “mathematical” derivative \(f'(x) = 1\).
- **identity**: The function \(f(x) = x\). This is an important comparison because it shows us what a “true” linear (both mathematically and computationally) network is capable of.

**Flattened models.** For the transfer functions that are mathematically linear, we can also compute the equivalent linear model. This just consists of a single dense layer, using the identity transfer function, that computes the linear function of the input. These appear in the results as “flat” variants.

**MNIST.** The first problem is the Modified National Institute of Standards and Technology handwriting dataset (MNIST). This is a standardized dataset of handwritten digits \((0-9)\) as \(28 \times 28\) greyscale images. This is chosen partly for trollish reasons. It dates from 1998, and even at the time of publication, accuracy with neural networks (98.4%) and other techniques (99.2%) were already extremely high [15].

For this problem, I augmented the dataset by randomly offsetting the training images by up to two pixels in any direction, and by adding Gaussian noise. The model’s input layer is just the \(28 \times 28\) greyscale values, and the output is a prediction for each of the ten digits. The models had two convolutional layers \((64 \times 3 \times 3)\) features, fully overlapping + 128 \(8 \times 8\) features, fully overlapping; then 32 \(128 \times 128\) features + 32 \(256 \times 2\) features with no overlap), then two sparse layers of 1024 nodes each, then a final dense output layer. The same initial weights and connectivity was used for each experiment. Internal layers use the transfer function being evaluated, but the output layer always used the identity transfer function. This is not a good choice for this problem (softmax makes more sense since the output is categorical) but I wanted the linear models to be truly linear. Using the same transfer function would have also disadvantaged functions with limited output range; downshift2 for example can technically output 1.0, but only for very large inputs (8192.0). The final identity layer can easily scale the useful range of the transfer function to the nominal range of the output. (This is essential for the chess problem below, where the output instead ranges \([-1, 1]\).)

See the source code for various hyperparameter settings (although if you are trying to learn good settings for hyperparameters, my code is not the place to look). I used the ADAM weight update trick [12], which does give me much better results than plain SGD in my experiments.

Results for MNIST are in Figure 4. A nice bug appears in Figure 5.

**CIFAR-10.** Another classic dataset comes from the Canadian Institute For Advanced Research. They capitalize “For” so that the acronym can be pronounced nicely. I mean to be fair MNIOSAT would have a certain ring to it too. This dataset contains 60,000 RGB images of size \(32 \times 32\), that are labeled into 10 different spirit animals: Airplanes, cars, birds, cats, deer, dogs, frogs, horses, ships, and trucks [14]. It is very similar to the handwriting prob-

<table>
<thead>
<tr>
<th>transfer function</th>
<th>flat accuracy</th>
</tr>
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<tbody>
<tr>
<td>logistic</td>
<td>98.20%</td>
</tr>
<tr>
<td>tanh</td>
<td>98.93%</td>
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<tr>
<td>leaky-relu</td>
<td>99.39%</td>
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<tr>
<td>plus64</td>
<td>82.66%</td>
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<tr>
<td>grad1</td>
<td>97.29%</td>
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<td>identity</td>
<td>81.96%</td>
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<td>downshift2</td>
<td>94.45%</td>
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<tr>
<td>plus64 ×</td>
<td>82.01%</td>
</tr>
<tr>
<td>grad1 ×</td>
<td>39.19%</td>
</tr>
<tr>
<td>identity ×</td>
<td>81.98%</td>
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9Learning with this function might work better if we instead approximate the derivative by something non-constant, like by computing the derivative of a smoothed version. However, due to implementation tricks in Tom7Flow, we need a derivative that is expressed in terms of the transfer function’s output (i.e. \(g(f(x)) = f'(x)\)); we would not be able to express the smoothed derivative because there are only 25 distinct values of \(f(x)\) in the \([-1, 1]\) range.
lem but more challenging (state of the art accuracy is more like 96.5%). You would struggle sometimes to figure out what these tiny thumbnails are, to be honest. Like with MNIST, I augmented the training set by randomly shifting the images and adding Gaussian noise. The network structure is the same as in the MNIST problem, except that in the first convolutional layer, each window is three times as wide to account for the three color channels.

Results for CIFAR-10 appear in Figure 6. One of the nice things about using standard problems is that we can understand how the results stack up against other researchers. Consulting a leaderboard of public results I see that the worst publicly known accuracy for CIFAR-10 is 75.86% [21]. The best result for the current work, using the sensible Leaky Relu transfer function, is 73.11%. So this is... last place. That’s actually pretty good; last place is the last winner (or the first winner, when counting from the end). Not to mention that we can get into even laster place by using the other exotic transfer functions. Even putting aside their aesthetic appeal, I feel that these inferior transfer functions are an important contribution to the field, as it seems to me that AI is getting too good, and too fast! Let’s take it easy there, guys!

**Chess.** This problem attempts to learn a good evaluation function for chess boards. Training examples are real chess positions (from the Lichess database) evaluated by a strong chess engine (Stockfish [30]). Stockfish generates two classes of scores: “Mate in N” if one side is known to have a series of N moves that wins (but “Mate in 1” is still better than “Mate in 4”), or a more subjective score, measured in pawns. (The score in pawns can seemingly be higher than 64, which is kind of funny because how are the pawns gonna fit on a 64-square board? DUAL-WIELD?) Mate is of course categorically better than the pawn score, as it is exact. Anyway, I squash this score into the range $[-1, 1]$ and that becomes the training instance. This network’s first layer has 256 $3 \times 3$ convolutional features, overlapping, as well as $32 \ 1 \times 1$ and $128 \ 8 \times 1$ and $1 \times 8$. Each of these is measured in terms of squares on the board, but each square actually corresponds to 13 inputs, for the 13 possible things that can be in that square (exactly one set to 1.0). We also have some non-square inputs, like the castling privileges and en passant state. So it’s not just the convolutional features but some sparse nodes too. And then we have some more layers (you can check out the source code if you really care about these details, which I doubt!) and then a final dense layer with a single output using the identity transfer function as before. No training

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<td>logistic</td>
<td></td>
<td>56.83%</td>
</tr>
<tr>
<td>tanh</td>
<td></td>
<td>67.82%</td>
</tr>
<tr>
<td>leaky-relu</td>
<td></td>
<td>73.11%</td>
</tr>
<tr>
<td>plus64</td>
<td></td>
<td>43.60%</td>
</tr>
<tr>
<td>grad1</td>
<td></td>
<td>53.56%</td>
</tr>
<tr>
<td>identity</td>
<td></td>
<td>41.07%</td>
</tr>
<tr>
<td>downshift2</td>
<td></td>
<td>46.54%</td>
</tr>
<tr>
<td>plus64 ×</td>
<td></td>
<td>32.76%</td>
</tr>
<tr>
<td>grad1 ×</td>
<td></td>
<td>30.58%</td>
</tr>
<tr>
<td>identity ×</td>
<td></td>
<td>41.04%</td>
</tr>
</tbody>
</table>

Here’s an idea for a SIGBOVIK paper: What’s the highest scoring chess position, according to Stockfish, for which it cannot deduce mate? One logistical challenge is that it seems to top out at +99, such as on this position (still no mate at depth 89).
data augmentation here (we have a basically limitless supply of positions to train on), but I do normalize the board so that it is always white to move.

For chess we can compute the accuracy, comparing to Stockfish as ground truth (Figure 7). We can also use the evaluation function to play chess. These chess “engines” just look at the possible legal moves and take the move that is most favorable, using the learned evaluation function (no game tree search). Playing against the best of these (“leaky”) it subjectively makes decent moves most of the time and can even beat me playing casually. I noticed that it had a lot of trouble “sealing the deal” in totally winning positions (which is not unusual for engines that don’t do game-tree search or use endgame tables), but the problem was actually more shallow: Due to a bug in the way training examples are gathered, the models were never exposed to checkmate or stalemate positions! Since training takes several days per function and the iron-fistedly punctilious SIGBOVIK deadlines were imminent, there simply wasn’t enough time to retrain them with access to these positions. However, since mate is a mechanical fact of the game (like what moves are legal) it seemed reasonable to fix this in the engine itself: When considering all the legal moves to make, it infinitely prefers a move that results in checkmate, and considers a move resulting in stalemate to have score 0.0, and otherwise uses the evaluation function. These “fix” versions of each engine perform very significantly better, although they likely overestimate the performance we’d get by actually fixing the model; there’s no guarantee that it would be able to accurately recognize mate, and the fixed versions’ greedy strategy of taking mate in 1 is always advantageous.

These players compete against each other as well as the engines from the Elo World project [26], giving a sense of their strength in an absolute scale (Figure 8). The raw versions perform reasonably; they all work better than a simple engine like “take the move that minimizes the number of moves the opponent will have,” \( \min \text{oppt moves} \). The fixed versions are much better, as expected. The “linear” engine using the grad1 transfer function, is competitive with the NES Chessmaster engine, and outperforms a 50% dilution of Stockfish. This is pretty solid given that it is doing no explicit game tree search. In fact (aside from the wrapper implementing the rules of chess and finding the maximum eval score), it is only performing a fixed expression of floating point addition and scaling! We could make this even more ideologically pure using techniques from Section 7.3.

**What transfer function is best?** The results on each of these problems are similar: The “leaky rectified” transfer function is best across the board. For data augmentation here (we have a basically limitless supply of positions to train on), but I do normalize the board so that it is always white to move.

For chess we can compute the accuracy, comparing to Stockfish as ground truth (Figure 7). We can also use the evaluation function to play chess. These chess “engines” just look at the possible legal moves and take the move that is most favorable, using the learned evaluation function (no game tree search). Playing against the best of these (“leaky”) it subjectively makes decent moves most of the time and can even beat me playing casually. I noticed that it had a lot of trouble “sealing the deal” in totally winning positions (which is not unusual for engines that don’t do game-tree search or use endgame tables), but the problem was actually more shallow: Due to a bug in the way training examples are gathered, the models were never exposed to checkmate or stalemate positions! Since training takes several days per function and the iron-fistedly punctilious SIGBOVIK deadlines were imminent, there simply wasn’t enough time to retrain them with access to these positions. However, since mate is a mechanical fact of the game (like what moves are legal) it seemed reasonable to fix this in the engine itself: When considering all the legal moves to make, it infinitely prefers a move that results in checkmate, and considers a move resulting in stalemate to have score 0.0, and otherwise uses the evaluation function. These “fix” versions of each engine perform very significantly better, although they likely overestimate the performance we’d get by actually fixing the model; there’s no guarantee that it would be able to accurately recognize mate, and the fixed versions’ greedy strategy of taking mate in 1 is always advantageous.

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**What transfer function is best?** The results on each of these problems are similar: The “leaky rectified” transfer function is best across the board.
fer function is generally best or close to best. The identity transfer function, which yields a simple linear model, is generally worst or close to worst. The sigmoid functions are all over the place. It is known that they are prone to vanishing gradients in deep networks, and I may simply have unfavorable hyperparameter settings for them. The experimental downshift2 function is generally bad, perhaps because its output is strictly positive or it has such a small dynamic range. Its shape also seems prone to the vanishing gradient problem. The small amount of nonlinearity introduced by plus64 does appear to give it a small edge over the identity, but its lack of an interesting derivative and the fact that it only produces a small number of output values are limiting. Importantly, the grad1 function—the centerpiece of the first third of this paper—performs decently on all problems. It clearly outperforms the linear models, despite being “linear.”

It is also interesting to compare the flattened versions of the linear transfer functions. These are the computed (mathematically) equivalent single-layer linear models. For plus64 the flattened version is worse in all cases; the unflattened model is taking advantage of the discretization in some way. For grad1 it is dramatically worse, both because grad1 models are substantially using the roundoff error and because the mathematical version of this function \( f(x) = x \times 1.036535 \) is not even a good linear approximation of the actual result (e.g. \( \text{grad1}(1) = 1 \)). Finally, the result for the identity transfer function should be mathematically equivalent, but it does not always produce the same results. This is unsurprising since we know that floating point calculations are not perfectly accurate, but it does hint that deep networks may make use of floating point roundoff internally, even if they are not using silly transfer functions!

Having proved the professor wrong, we could stop there, but did huge mathematical breakthroughs ever arise from taking the option to stop there?!

5 Non-monotonic functions

Because of the way that addition and scaling are defined (do the real mathematical operation, then round), they preserve monotonicity: If \( x \geq y \), then \( f(x) \geq f(y) \). But this is only true if we limit the form of the function to a series of additions of constants and (non-negative) scaling. There are other expressions that are mathematically linear but don’t take that form; for example:

\[
\begin{align*}
  f(x) &= x - 4096 - x + 4096 \\
\end{align*}
\]

This is of course mathematically equivalent to \( f(x) = 0.0 \), but with half precision it is a square wave function (here pictured \([-8, 8]\)):

For some values of \( x \) the terms cancel out, and for others the rounding error compounds. This function is not as well-behaved as it appears: the first pulse has width 0.99609375 and the second has width 1.

Here is \( f(x) = \text{grad1}(x) - x \), which is also linear:

Generally speaking, we can create a large variety of functions by computing the interference patterns between other functions, since the sum or difference of two “linear” functions is also “linear.” In general we’ll consider expressions of this form:

\[
E \ ::= \ x \\
| \ E \times \ c \\
| \ E + \ c \\
| \ E + E
\]

Where \( x \) is the function variable, and \( c \) is one of the 63,488 finite half-precision constants. We can derive negation \( (E \times -1) \) and subtraction of constants \( (E + -c) \) and expressions \( (E + (E \times -1)) \) since every number has an exact negation by flipping its sign bit. Exact division is possible when \( 1/c \) is representable, and there is almost always a close approximation.

This formulation leads to a tempting approach for approximating a function iteratively, like a Taylor series. Given a target function like \( \sin(x) \), we can begin with an approximate expression for it, like \( x \), and then add and subtract terms to improve the approximation. I don’t know of any systematic way to improve the approximation at each step (they are not well-behaved mathematically, and I am not good at math), but by using computer search I can sure make some complicated functions with many different shapes.

An approximation of \( \sin \) appears in Figure \( \text{fig:sin} \). It is fun to watch an animation of the successively improving approximations, but you can’t see that since you’re reading an old-fashioned paper. Perhaps you can find a video of this at tom7.org/grad.

5.1 Fractals

Next, I endeavored to deploy these functions for something useful: Fractals. Famously, fractals are simple functions with complex (often literally) behavior. For example, the Mandelbrot set considers each complex point \( c \) (plotted on the plane as \( x + yi \)) and computes whether \( z_i = z_{i-1}^2 + c \) diverges or not. It’s lovely, but squaring is not linear!
What if we just create a linear function that approximates \( f(x) = x^2 \)? This is definitely possible, using the approach described above. After 184 successful error-reducing rounds we get the following approximation, with 112,204 linear operations:

Aside from the funny business near the origin, this is a fairly accurate approximation of the square function, so you might hope that it would draw a perverted Mandelbrot set. Unfortunately, it produces a much sadder blotch (Figure 10). To see why, consider the normal definition of squaring for a complex number:

\[
(a + bi)^2 = a^2 + 2abi + b^2i^2 = a^2 + 2abi - b^2
\]

Note that the real coefficient \( a \) ends up part of the imaginary coefficient \( 2ab \) in the result, and the imaginary coefficient \( b \) becomes part of the real part (because \( i^2 \) is real). This means that squaring a complex number cross-pollinates between the two components, yielding a kind of wacky rotation if we think of them as 2D coordinates.

But here, squaring is approximated as a series of operations of the form \( w_1 + w_2 \) and \( w \times c \) for constants \( c \). These operations on complex numbers are less interesting:

\[
(a_1 + b_1i) + (a_2 + b_2i) = (a_1 + a_2) + (b_1 + b_2)i
\]
\[
(a + bi) \times c = ac + bc \text{i}
\]

Alas, these operations are boring; the real parts always stay real and the imaginary parts always stay imaginary. This is why the crummy blotch has all sorts of vertical and horizontal lines in it: As we iterate the function we are iterating two independent components, and the resulting picture is just some interference pattern between them.

This seems pretty definitive. Even if we had some kind of hardware implementation of complex numbers with rounding error to abuse, there would be no reason to have the linear operations do any cross-pollinated rounding. Professors take note: The complex numbers do provide some refuge!

Still, a lot of chaos can emerge from these functions that should not be possible with “linear” ones. For example, here is a complicated function made by stringing 36,637 addition and scaling operations together:
Iterating this function produces chaotic results because of its nonmonotonicity. In Figure 11 I plot (using color) the magnitude of $z$ after 256 iterations of

$$z_i = f(z_{i-1}) \times c$$

This is mathematically linear (as $c$ is a constant and $f$ a linear function). Nonetheless, it produces an appealing picture. I think this is a fractal in the sense that it is chaotic, has a color gradient, and could be on the cover of an electronic music album. It is not a fractal in the sense that if you zoom in on it, you get infinite detail of self-similar shapes. In fact, if you zoom in on it only a modest amount, you encounter rectangular pixels as you reach the limits of half-precision floating point. (And because this fractal is built by abusing those very limits, it is not even possible to get more detail by increasing the accuracy!)

5.2 Bonus digression: Baffling numbers

Imagine you are my professor. You assign a class project to “make fractals using floating point roundoff error,” for some reason. You spot me in the computer lab and I’m obviously way off track, because on-screen is some kind of 3D fractal. The Mandelbrot set cannot be extended to three dimensions, you say, because of the Frobenius theorem: Only algebras of dimension 1 (real numbers), 2 (complex numbers) and 4 (quaternions) work [8]. Unclear how the professor speaks the citation aloud in this scenario. I say I “know” this fact, but I “don’t care.” You say that my three-dimensional algebra can’t be associative, because that’s “just a mathematical fact.” I say you know what else isn’t associative? The floating point numbers, my dude.

Enter the baffling numbers, ill-advised extensions of the complex numbers to three dimensions. Here we have numbers of the form $a + bi + cj$. Addition is just pointwise, and there are several options to complete the story for multiplication, namely the values of the cells $A$, $B$, and $C$ in this table:

<table>
<thead>
<tr>
<th></th>
<th>1</th>
<th>i</th>
<th>j</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>1</td>
<td>i</td>
<td>j</td>
</tr>
<tr>
<td>i</td>
<td>i</td>
<td>−1</td>
<td>U</td>
</tr>
<tr>
<td>j</td>
<td>j</td>
<td>V</td>
<td>W</td>
</tr>
</tbody>
</table>

The cells $U$, $V$, and $W$ are baffling numbers (i.e. each some $a + bi + cj$). Some choices are degenerate, but this gives us a family of options. It is known that no matter the choices, this does “not work” (in the sense that the resulting algebra is not associative) but we don’t need associativity to draw fractals. Plus, who’s gonna stop me, FROBENIUS?!

I tried a few options, but thought that $U = i$, $V = j$ and $W = 1$ produced satisfyingly trippy results. The Mandelbrot is straightforwardly generalized to the “Bafflebrot” (the starting point $c$ is just a baffling number now; everything else is the same). I generated a 3D object by defining an implicit surface based on whether a sampled point is still inside the set after 25 iterations, using Marching Cubes [17] to discretize it. The resulting mesh is 2 gigabytes and crashes every library that attempts to programmatically simplify it. I do admire and encourage its defiant spirit. A rendering appears in Figure 12.

Drawing fractals is fun and everything, but I grew weary of the exercise because there is no real goal other than to make a cool picture. Instead I turned to something with a clearer challenge to overcome: Linear Cryptography.

6 Linear cryptography

Cryptography is like fractals minus drugs. One of the most basic components of cryptography is a pseudorandom num-

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8Or else is equivalent to the complex numbers.
Figure 12: The 3D “bafflebrot” sliced in half and projected to 2D. This fractal was created with the “illegal” number system called the baffling numbers. They’re like the complex numbers but more so. The object is truncated along its j axis, showing a perfect ripe Mandelbrot inside.

Figure 13: The substitution-permutation network that forms a half decent pseudorandom number generator. The same substitution (“s-box”) is applied to each byte. Then the 64 bits are permuted. Finally, bytes are combined with modular addition and subtraction. This function passes the “Big Crush” suite and can be implemented with only half-precision floating point addition and scaling.

A B C D E F G H

S-Box S-Box S-Box S-Box S-Box S-Box S-Box S-Box

The basis of this function is the classic substitution-

a symmetric encryption algorithm like AES, with the key hidden, has this property. Practically speaking, though, we can subject the function to a wide variety of statistical tests, and if it looks random to every test, then this gives us good confidence.

Specifically, my goal is to design an algorithm that takes 64 bits of data (represented as half-precision floats) to another 64 bits, such that the stream of low-order bits from iterating this function passes the TestU01’s “Big Crush” suite of 106 statistical tests. This suite is a successor to Marsaglia’s “DieHard” battery of tests, itself an improvement on Knuth’s tests from The Art Of Computer Programming.

The basis of this function is the classic substitution-


12 Truly good cryptographic algorithms are also openly studied by experts. Of course nothing in here is to be used seriously, and not just because these algorithms are ridiculously slow. But I guess if you are stuck on a desert island with only the floating point addition and scaling operations, and a copy of this paper, then it would be a reasonable starting point for encrypting your messages. I do not recommend, if stranded on a desert island, to send encrypted messages: They may not be readable to your potential rescuers!
permutation network. First, each of the eight bytes are substituted with a different byte using a table (this is the mathematically non-linear step). Then, the 64 bits are permuted. Finally, some of the bytes are modified additively. An illustration appears in Figure 13.

The substitution table (“s-boxes”) was generated by computer search with an objective to maximize the “avalanche property” (when a bit of the input is complemented, about half of the output bits should be complemented). The permutation was generated to maximize dispersion; each quartet sends each bit to a distinct quartet in the output. This is not the important part. We could have just used known good tables.

To implement this with half-precision floating point, we could represent each bit with its own half, but that is no fun. The state will be represented with eight half-precision floats, each representing one byte’s worth of information. Since we have been fixated on the $[-1,1]$ interval so far, a byte will be stored as any value in $[-1,1]$. This is not the important part. We could have used $[−128, 128]$ instead. The one just pictured is one of the smallest expressions; speaking of which, my first approach was to try to approximate these expressions, but I discovered them with computer search and some elbow grease. We’ll call this one Choppy functions. I’ll say now that the value easily gets stuck in short cycles due to this inaccuracy.

I spent a long time writing code to simplify these expressions and generate LaTeX for them, by the way! As usual, I thought it would look cool when I got it working, but it just looks like a bunch of numbers.

We can think of this function as a basis vector, representing the 256-dimension vector $(0, 0, 0, \ldots, 0, 1, 0, 0, 0, 0, 0, 0)$. We’ll call this one $b_{249}$ since it selects the integer 249. If we can find $b_n$ for each $n \in \mathbb{Z}_{256}$, then we will be able to combine them to systematically construct functions.

The one just pictured is one of the smallest expressions; most are much larger. I wish I could tell you that I figured out the principles underlying how to analytically generate these functions, but I discovered them with computer search and some elbow grease.

The function produces the same result for any representation of an integer, and that result is the smallest representation of an integer. These functions are maximally useful in that they are “liberal in what they accept,” but most are much larger. I wish I could tell you that I figured out the principles underlying how to analytically generate these functions, but I discovered them with computer search and some elbow grease.
but “conservative in what they return” [28]. It also means that each function also can be understood as a function $\mathbb{Z}_{256} \to \mathbb{Z}_{256}$, so we can represent them as a vector of 256 integers. The basis vectors $b_n$ are those that are of the form $(0, \ldots, 1, \ldots, 0)$.

I then conducted computer search for choppy functions, putting those into a database (keyed by the corresponding integer vector). Some are easy to find, others harder. Summing and scaling choppy functions yield choppy functions (as long as the vectors remain integral and in range), so I use a simplified version of Gauss-Jordan elimination [32] to solve for basis vectors. Once I have $b_n$, this column can be changed at will for any existing choppy function (by just adding or subtracting multiples of $b_n$), so new choppy functions that only vary in that column can be ignored.

By trying a variety of operations that are known to be useful (e.g. iterated multiplication of constants near 1.0) and hill-climbing towards functions with the choppy property, it is not too hard to find $b_n$ for most $n$. It seems to become more challenging for $n$ near 128; this is the point 0.0 in half-precision. Specifically, the hardest problem was to make a function that produced different results for inputs $< 0.0$ versus inputs $\geq 0.0$. This is the zero-threshold problem.

**Why is this hard?** Distinguishing between negative and non-negative numbers is deceptively difficult. Looking back to the function $f(x) = x + 128.0 - 128.0$ (Section 4), it has useful discontinuous steps, but note that the discontinuity does not happen at zero. This is because we are rounding to the nearest value, and so small negative numbers near zero end up rounding to the same result that zero does. Moreover, the resolution of the floating point numbers is highest near zero (especially because of subnormal numbers), which exacerbates our attempts to control rounding of them. For example, you might think that we could simply shift this function left and right by substituting $x + c$ for $x$ in its body. This would work mathematically, but it does not work for floating point numbers, because each operation performs some rounding. If this rounding ever ends up conflating a negative number with a non-negative one, we will not be able to recover.

I found a zero-threshold function using a combination of manual and computer search. This was some ordeal, and the resulting enormous function is in Figure 14. Perhaps you are smarter than me and can find a better one!

**Substituting and permuting.** In any case, with this function it was possible to form a complete basis. This basis makes it “easy” to perform operations on half values that represent bytes. For example, the s-box step substitutes some distinct byte for each different input byte. This would normally be implemented with a table lookup. If we compute $b_n(x) \times \text{subst}[n]$, this returns the correct result $\text{subst}[n]$ if the input $x = n$, and 0 otherwise. So if we just sum all 256 of these up, exactly one of them will be nonzero, and the correct substituted value.

Permutation is defined on the component bits. Here, we compose a function that computes each of the eight output bytes. We use the same approach of summing a bunch of $b_n(x)$ evaluations (each multiplied by the correct answer).

Here we are testing whether the input has some particular bit set (a sum of the 128 $b_n(x)$ functions where $n$ has that bit set), and the output is the power of two that sets the appropriate output bit. Many of these functions would have simpler implementations (for example, “is the high-order bit set?” is the same as the zero-threshold function) but at this point I was happy to just have something working, and taking some joy in how absurdly large the functions were getting.

The cipher also includes addition and subtraction mod 256. Addition and subtraction are already available for half-precision floats, and they have faithful behavior, so we just need to implement the wrapping-around behavior so that the result is strictly in $[-1, 1]$. This is straightforward with the zero threshold function [28] we produce corrective factors if the result is $\geq 1$ or $< -1$ (zeros otherwise). We then add those corrective factors produce the remainder we desire.

### 6.1 Benchmark results

To evaluate the quality of the pseudorandom number generator, I used the TestU01 “Big Crush” suite. This test needs a sample of 1.64 billion bits, so I actually evaluated it on equivalent code that performs the steps using normal integer operations. Even then, the suite takes several days to run, so I modified it to run tests in parallel and cache the results of completed tests. This saved me from losing data if my computer crashed or needed to be rebooted.

Results appear in Figure 15. Passing these tests does not ensure that the pseudorandom number generator is good for cryptography, although it is a good start.

Running single-threaded on a 3.0 GHz Threadripper 2990WX, this function generates 25.8 bytes of randomness per second, which is slow. By precomputing the substitution, permutation, and zero threshold expressions (so they can be performed by lookup into 64k-entry tables), it generates 18,685.2 bytes per second, which is still slow.

If we were building an encryption algorithm (a symmetric block cipher), it would be natural to use this as its “round function.” In a Feistel network [7], each input block (128 bits) is broken into two halves; one of them is mixed with some key bits (for example with XOR) and then passed to this function. Its output is XORed with the other half; the two halves are swapped, and this “round” is repeated many times until we believe that the data are suitably screwed up. Decryption is the reverse. We can use addition and subtraction mod 2^8 to combine the data instead of XOR

14Technically we need to do some multiplicative adjustments to put the value in $[-1, 1]$.

15Compare to the remarks “why is this hard?” above. Here, $zt(x - 1)$ does do what you’d want, shifting the threshold value from 0 to 1. This is because there is less precision near one than near zero.
7 THE ULTIMATE THROW-BACK

Having developed a basis for extracting arbitrary bits, we can express any function of a single variable, and we’ve seen how some other functions (like addition mod 2^8) can be done. At this point, it seems like we probably have the building blocks to demonstrate that addition and scaling on half-precision floats is Turing complete. I mean, pretty much everything is Turing complete. In the past, I built computers that were perfect and beautiful, such as a hardware implementation of the NaNDy 1000, a computer architecture that computes using only floating point NaN and Infinity [27]. In a concession to ideological purity, though, the NaNDy 1000 has no I/O. So it is very boring to use.

For today’s investigations of the capabilities of floating point, I’ll make the opposite concession: Let’s make a computer that is exciting to use, but that makes some (reasonable) ideological concessions so that it can do something interesting.

7.1 fluint8

First of all, if we want to do some serious computation, 25.8 bytes per second isn’t going to cut it. To look for performance enhancing substances, I perused the back catalog of the world’s most prestigious conference, SIGBOVIK. There in the 2018 edition, on page 125, I found an intriguing paper, The fluint8 Software Integer Library, by Drs. Jim McCann and . . . Tom Murphy VII? Wait, that’s me? I already wrote this paper?! [20]

The fluint8 library represents an element of \( \mathbb{Z}_{256} \) (a.k.a. uint8) as a 32-bit float, and provides multiplication, addition, subtraction, negation, division, and bitwise functions and, or, and exclusive or.

Compared to the approach discussed in Section 6 using “choppy functions,” fluint8 has much more simple and sensible implementations of functions like addition:

```c
inline float fu8_add (float a, float b) {
    float x = a + b;
    x -= x - 127.5f + 3221225472.0f - 3221225472.0f;
    return x;
}
```

The \( x \rightarrow x \ldots \) line applies the corrective factor to implement wrap-around, which we previously did using the zero-threshold function. Why can it be done so much more simply here? First, fluint8 represents \( n \in \mathbb{Z}_{256} \) as \( n \), so a number like 27 is represented as 27.0 instead of, say, \(-1 + 2^7/128\). Second, it requires that the number be represented exactly as this value. The figures in the fluint8 paper are somewhat misleading as they are plotted only

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Figure 15: Results of the TestU01 “Big Crush” suite on the pseudorandom number generator built from floating point roundoff error. A p-value of < 0.001 or > 0.999 is considered suspect by the suite, so all tests pass here.
for input values that are already exact integers; if we test \texttt{fu8\_add} on values like 100.1875 and 11.0703125 we do not get 111 (Figure 16). On the other hand, this is a very reasonable choice to make; we can simply have a representation invariant that only one of these 256 values is used, and preserve that invariant with every operation. It won’t work great for the continuous domain (e.g. plotting fractals) but is a much more practical choice for discrete data (e.g. encryption). Since I like to work at the intersection of Theory, Impractice, and Practice, this is appealing!

But: The library uses several operations that are not linear! In particular, its implementation of bitwise functions like XOR perform squaring and multiplication of the two arguments. It was not a design goal of \texttt{fluint8} to use only addition and scaling, but it is a design goal today, so we must address that.

### 7.2 hfluint8

The use of nonlinear operations is a problem we will rectify, forthwith, but the other ideas are suitable for building a computer. In the \texttt{hfluint8} (for half float linear unsigned int 8-bit) library, a \texttt{hfluint8} will be represented by a single half-precision floating point number, and always one of the exact integral values in \([0, 256)\).\(^\text{17}\)

```
struct hfluint8 {
    half h;
    // ...
};
```

Let’s begin with one helper function.\(^\text{18}\)

```
half RightShiftHalf8 (half xh) {
    half SCALE = GetHalf (0x1c00); // 1/256
    half OFFSET1 = GetHalf (0xb7f6);
    half OFFSET2 = GetHalf (0x66b0);
    return xh \* SCALE + OFFSET1 + OFFSET2 - OFFSET2;
}
```

If the function is given an integral half \(xh\) in \([0, 512)\), it returns \(xh \gg 8\). This value is always 1 or 0. The calls to \texttt{GetHalf} interpret a 16-bit constant as a half, which is useful to be precise (many decimal expressions like 0.1 are not exactly representable in floating point). I also found that if you use literals like 0.00390625_h, the code runs much much more slowly because it inhibits some optimizations or perhaps the user-defined literals are parsed at runtime (?). Aside from wanting to avoid operations like parsing that might not be addition and scaling on halves, we will struggle with performance of these functions as we use them for real

---

\textsuperscript{17}In fact, all integers from -2048 to 2048 are available, so we could consider implementing signed 11-bit numbers in a future \texttt{hfsint11} library.

\textsuperscript{18}These code samples have been simplified to fit the extremely capricious SIGBOVIK column width requirements. For example, \texttt{GetHalf} is a \texttt{constexpr} function, so these constants are really declared as \texttt{static constexpr} and computed completely at compile time. See the full code and verify that it complies with the rules at \url{https://sourceforge.net/p/tom7misc/svn/HEAD/tree/trunk/grad/}.

Figure 16: \textbf{Top:} Error of the \texttt{fluint8} addition function on general floating point values in \([0, 256)\). This is a detailed zoom of the region \(x \in [252, 256)\) and \(y \in [0, 4)\), but the rest of the image is almost identical. Each pixel compares the \texttt{fluint8} sum of \(x\) and \(y\) to the expected value \((\lfloor x \rfloor + \lfloor y \rfloor \mod 256)\). The top-left pixel in each cell is the case where \(x\) and \(y\) are integers; we get the correct result (no error). All other pixels are wrong, either too high (green) or too low (red).

\textbf{Bottom:} Same with the error of the \texttt{floor} of \texttt{fluint8}’s sum function. This shows that the output is usually not even in the correct interval. However, observe the multitude of Triforces!

\textbf{Nowhere, or a lot of places if you think about it:} The modular addition operation from Section \texttt{6} is not pictured for comparison because it would be all white, meaning no error. You can actually imagine it occupies any blank portion of this paper, such as the inner hole of a letter ‘o,’ or the entire back of a page if printed single-sided.

Graphics produced using \texttt{ImageRGBA} computational for loop engine.
computing. Anyway, we are just dividing by 256 (by multiplying by $1/256$) and then adding some mysterious constants to ensure that the result is exactly 1 or 0.

Next, we can perform addition:

```c
hfluint8 hfluint8::Plus(hfluint8 x, hfluint8 y) {
    half HALF256 = GetHalf(0x5c00); // 256.0
    half z = x.h + y.h;
    half o = RightShiftHalf8(z);
    return hfluint8(z - o * HALF256);
}
```

As in `fluint8` we can simply add the arguments, giving a result in $[0, 512]$. The shift function just discussed then allows us to compute 1 if the value is out of range or 0 otherwise. We multiply this by a corrective constant (256.0) and subtract that away. So easy.

For all other operations we work on the domain $[0, 256]$. We also have a right shift by one bit:

```c
half RightShiftHalf1(half xh) {
    half SCALE = GetHalf(0x37fa); // 0.4986...
    half OFFSET = GetHalf(0x66cd); // 1741.0
    return xh * SCALE + OFFSET - OFFSET;
}
```

Right shifting is integer division by two. Roughly we are dividing by two and then offsetting to a part of the floating point number line where only integers are representable, then offsetting back. However, with a constant of exactly 0.5 some of the rounding would be in the wrong direction; the constant 0.49853515625 just happens to work.

We can shift by multiple places by repeating this operation multiple times. However, the library has direct solutions for several other shift distances, since this is more efficient than repeating a single shift.

Next, bitwise operations. These are all based on the AND function:

```c
half BitwiseAndHalf(hfluint8 a, hfluint8 b) {
    half result = GetHalf(0x0000);
    for (int bit_idx = 0; bit_idx < 8; bit_idx++) {
        // Low order bit as a - ((a >> 1) << 1)
        hfluint8 as = RightShift1(a);
        hfluint8 bs = RightShift1(b);
        half a_bit = a.h - LeftShift1Under128(as).h;
        half b_bit = b.h - LeftShift1Under128(bs).h;
        // Computes 2-bit_idx. A constant.
        half scale = GetHalf(0x3c00 * 0x400 + bit_idx);
        half and_bits = RightShiftHalf1(a_bit + b_bit);
        result += scale * and_bits;
    }
    return result;
}
```

This function shifts each input down 8 times, stripping off the low order bit at each step. Note that since we run this loop exactly 8 times, it can simply be unrolled, removing any whiff of non-linearity, and the constants computed at compile time. `LeftShift1Under128(x)` is just $x + x$ without any need to worry about modular arithmetic, as it cannot overflow.

An interesting line is the computation of `and_bits`, which is the logical AND of the low-order bit from $a$ and $b$. In `fluint8` we simply compute $a_{\text{bit}} * b_{\text{bit}}$. This has the correct value, but is not linear (observe that if we were to compute $x \& x$ we would end up squaring a function of $x$ here). Instead we compute $(a_{\text{bit}} + b_{\text{bit}}) >> 1$, which produces the correct result.

Being able to compute the bits in common allows us to easily derive OR and XOR:

```c
hfluint8 BitwiseOr(hfluint8 a, hfluint8 b) {
    half common = BitwiseAndHalf(a, b);
    return hfluint8((a.h - common) + b.h);
}
```

```c
hfluint8 BitwiseXor(hfluint8 a, hfluint8 b) {
    half common = BitwiseAndHalf(a, b);
    return hfluint8((a.h - common) + (b.h - common));
}
```

These subtractions and additions cannot overflow.

It will be common to perform bitwise operations with constants, so `hfluint8` supports versions with a compile-time constant argument, which can skip a bunch of work. These run about $5 \times$ faster.

We also have some operations that are not supported by `fluint8` but that we will need for the current project. A basic operation is to test for zero. `IsZero` returns 1 if the input is 0, or returns 0 for any other argument:

```c
hfluint8 IsZero(hfluint8 a) {
    half H255 = GetHalf(0x5bf8); // 255.0
    half H1 = GetHalf(0x3c00); // 1.0
    half nota = (H255 - a.h);
    return hfluint8(RightShiftHalf8(nota + H1));
}
```

For an input of zero, complementing it yields 255, and adding 1 overflows to set the 8th bit. So we shift that bit to the ones place and are done:

```c
half H255 = GetHalf(0x5bf8); // 255.0
hfluint8 mask = hfluint8((cc.h * H255) + (cc.h + H1));
return BitwiseAnd(mask, t);
```

This computes either the mask 00000000 or 11111111 and uses the existing bitwise AND operation. Bitwise AND is not fast, and it does more work than it needs to in this case because we know one of the arguments is all zeroes or all ones. It is faster to inline the bitwise AND routine but keep checking the ones place. Even better is this wild ride:

```c
if (cc, t) returns t if cc is exactly 1, returns 0 if cc is 0, and is otherwise undefined. A simple implementation of this is:

```c
half H255 = GetHalf(0x5bf8); // 255.0
hfluint8 mask = hfluint8((cc.h * H255);
return BitwiseAnd(mask, t);
```

This computes either the mask 00000000 or 11111111 and uses the existing bitwise AND operation. Bitwise AND is not fast, and it does more work than it needs to in this case because we know one of the arguments is all zeroes or all ones. It is faster to inline the bitwise AND routine but keep checking the ones place. Even better is this wild ride:

19Earlier iterations of this function were much more complex! For example, $x + 15 + 65248 - 65248 \times 0.03125$ maps 0 to 0, but any other number to some number in $[1, 15]$, and then a similar function compresses that range down to exactly 0 or 1. But sometimes you miss the obvious stuff until you start writing a paper about it for a prestigious conference. No doubt some other functions in here could be improved!
hfluint8 If(hfluint8 cc, hfluint8 t) {
    static std::array<half, 8> OFF = {
        GetHalf(0x77f9), GetHalf(0x7829),
        GetHalf(0x77fb), GetHalf(0x78e2),
        GetHalf(0x77fd), GetHalf(0x78b0),
        GetHalf(0x77ff), GetHalf(0x7864),
    };

    half HALF1 = GetHalf(0x3c00); // 1
    half HALF128 = GetHalf(0x5800); // 128
    half HALFNEG1 = GetHalf(0xbc00); // -1
    half HALF0 = GetHalf(0x0000); // 0

    half xh = t.h;
    half nth = HALF1 - cc.h;
    half c128 = HALF128 * nth;

    std::array<half, 8> COFF;
    for (int i = 0; i < 8; i++)
        COFF[i] = OFF[i] * nth;

    for (const half &h : COFF) xh = xh + h - h;
    for (const half &h : COFF) xh = xh + h - h;

    return hfluint8(xh * HALFNEG1 + HALF0);
}

The 8 constants in OFF, when added to and subtracted from a hfluint8, will always round such that the low six bits become 0. To have behavior conditional on cc, first we multiply each constant by 1-cc. This results in either the original constant or 0. If zero, then adding and subtracting them does nothing. Then we add and subtract those results, clearing the low six bits, and (conditionally, using the same trick of multiplying by the condition) subtract from 128. This clears the top two bits for the range of possible values (but may reset low-order bits). Then we add and subtract the sequence again, clearing the low six bits again. At the end we apply a corrective negation and then add 0 to avoid outputting 0 and we’re done.

hfluint16. Several other operations are available for hfluint8, like AddWithCarry, but we shan’t elaborate them all here, lest we construct hfluenza. One more concept is needed before we get to the application: 16-bit integers. The hfluint16 type is implemented as a pair of hfluint8 bytes. We will only need a small number of operations: Addition, subtraction, bitwise operations, sign extension of hfluint8 If, and stuff like that. These are all cleanly implemented in terms of the fluint8 operations like AddWithCarry.

### 7.3 Linear gameplay

Now we can build an 8-bit computer. I like to work at the intersection of theory, impractical, practice, and entertainment, and the most entertaining 8-bit computer is the Nintendo Entertainment System, so let’s build that. The full NES has many components (video output, controllers, sound, RAM, cartridge mappers), and it’s not even clear what it would mean to implement “linear” versions of these. So for this project we will replace the CPU, which is a variant of the Motorola 6502 called the Ricoh 2A03. Each instruction that the CPU executes will be done entirely with linear half-precision floating point operations. This is done in software emulation, upgrading a version of the FCEUX Emulator [6] that I forked many years ago [23].

The 2A03 has 8-bit registers A, X, Y, a stack pointer S and processor flags P. Each is represented as a hfluint8, of course. It also has a 16-bit program counter PC, which we represent as a hfluint16. Putting aside the many complexities, at each step it reads the byte at the program counter, which denotes one of its 256 instructions. It then executes the corresponding instruction, which produces new values for the registers and advances the program counter a variable amount. For example, a very simple instruction is TAX (0xAA), which could be implemented like this:

```
reg_X = reg_A;
reg_P = (Z_FLAG8 | N_FLAG8) & reg_P;
```

It is not implemented like this. Everything gets more complicated. But anyway, the TAX instruction Transfers (copies) the A register to the X register, and then updates the Zero and Negative bits of the flags register. We have all of these operations on hfluint8, so it’s just a matter of doing it.

**Memory.** For instructions that act solely on registers, this approach suffices. Most instructions read from or write to memory, including just to read additional arguments to the instruction. This is a problem because we don’t have any kind of branching; we always need to execute the exact same sequence of additions and scaling operations. We can work with this by computing condition codes: “Is this write actually happening, or are we just computing it because we always have to do the same sequence of operations?” Then a write `addr = val` can be made conditional using our If operation, like

```
mem_addr = If(cc, val) + If(1-cc, mem_addr)
```

This has other problems (for example when the address is not know at compile time, which is typical) but the biggest one is that all memory accesses on the NES are potentially effectful. This is because various things are attached to the memory controller that perform actions when addresses are accessed. For example, writing two consecutive bytes to 0x2006 will load them as an address into another chip (the PPU) and then writing to 0x2007 will write bytes into video memory at that address. Writing to 0x4014 will begin a DMA loop that copies 256 bytes from the main address space to video RAM, suspending the CPU for 512+ cycles. Reads can have effects as well, and these effects are not from a small set because they can include arbitrary hardware in the cartridge itself [25].

So here we have a sort of concession: We introduce two primitive operations `ReadIf(cc, addr)` and `WriteIf(cc, addr, val)`. These take a hfluint8 condition cc (exactly 0 or 1), a hfluint16 address, and (for writes) a hfluint8 value to write. If the condition code is 0, nothing happens, and an arbitrary value is returned. If 1, the read
or write takes place, including its side-effects. This would be a realistic model if we implemented a hardware version of this chip, which only used floating point operations internally; its hardware pins for interfacing with memory would simply include a bit for whether we actually want the read or write to happen. (The actual 2A03 pinout has a “R/W” pin, for example.)

Doing it correctly. The remainder is reasonably straightforward given the tools we’ve already built. One challenge is simply not screwing up. 256 instructions is a lot, and the original code is extremely awful; it is filled with macro hacks that assume specific variable names and values of constants, pirate jokes, references to mysterious global variables named stuff like temp, feuds between developers commenting out each other’s “wrong” code, and so on. As I developed the hfluint8-based emulator, I strove to keep the emulator in a working state as often as possible so that I could test it against the reference implementation. One technique was to do various pieces of code in easy, cheating ways, but to record each time I cheated by incrementing a global counter. Each time I replaced reasonable fast code with ideologically pure, non-cheating code, which is typically much slower, the cheating went down and the runtime went up; see Figure [17] This makes it like a game.

Another challenge is that the 2A03 has dozens of undocumented instructions with mysterious behavior. Most of these are not used by any game in my test suite, which means I run the risk of breaking one of these instructions and not knowing. Some of these instructions are very weird, since they are essentially the consequence of 6502 sub-units (designed for implementing other instructions) being connected together in ways that are not motivated by useful behavior. For example, the XAA instruction (0x8B) bitwise-ORs the A register with 0xEE (setting all but two bits), then ANDs with the X register, then ANDs with an immediate byte. Others are just as weird but much more complex. Since I want the emulator to be as complete and correct as possible, I wrote a new “game” that I could use as an additional test ROM (Figure [18]). This “game” executes dozens of undocumented instructions at startup, writing interesting state to RAM to create a record of their behavior. The game then displays the first half of RAM on screen. This gives some amount of protection against regression on these instructions.

Everything, everywhere, all at once. Each instruction is otherwise straightforward to implement. The remaining challenge has to do with the instruction dispatch. A natural way to write the instruction loop is to do switch on the instruction byte, but that is not a linear operation. Instead, we always execute all of the instructions. Before this, we make 256 copies of the CPU state (the registers); this is linear because it’s just copying a finite number of variables. Each copy also has an active flag (a hfluint8 with 1 or 0). We set this for exactly one of these instructions, by computing If(Eq(insn_byte, n)) for each of the 256 n. Then we execute each instruction on its copy of the state; it does all its computation, and any read or write is additionally conditioned on its active flag. This way only the active instruction’s memory accesses actually occur.

We then need to select the instruction that was actually executed and copy its state back to the “real” CPU state. We do this by conditionally clearing each register:

\[
\text{reg} = \text{If} (\text{active}, \text{reg})
\]

We then set the real CPU’s register to the sum of all of the registers from the instruction-specific states. Exactly one (the active one) will be nonzero, so we get that value. We use this same technique to keep track of how many cycles have elapsed, since various emulator timing depends on this.

A bad thing about this approach is that it’s more than 256 times slower than just executing a single instruction, and this is the main reason why the emulator is so slow. A good thing is that there is no cheating. Another good thing is that the instructions are all reading and writing distinct data, so they can actually be executed in parallel. The final benchmarks here are from running on 8 cores.

7.3.1 It’s a-fine, Mario!
The emulator can play any NES game supported by FCEUX (which is basically all of them; despite the horrors in this emulator’s code, it has great compatibility). My

![Figure 17: During the development of the emulator, the FPS achieved (blue) versus the number of times the code “cheats” due to incomplete implementation (red). Log scale. Honestly there’s not much to get from this except that we start with a lot of FPS (3500) and a lot of cheats (65 million) and end with few FPS (0.1) and no cheats. I guess it also shows that this took many iterations to implement. The reason that cheats does not monotonically decrease is that a single cheat (e.g. a switch on the instruction byte) can mask the need for hundreds of other cheats.](image-url)
benchmark was the first level of the classic *Super Mario Bros.*, playing sequence of 2210 inputs that completes level 1-1 in 36 seconds. The emulator runs this as fast (or as slow) as it can. Normal frame rate is 60 FPS. The original implementation runs at 3500 FPS; after many performance tweaks I got my *hfluint8* version to run at

0.1154 FPS

In print, the frame-rate is always zero, anyway (Figure 19). 8.6 seconds per frame is firmly in “not playable” territory, but it is tolerable for installation artwork, let’s say. I have played AAA titles that, at launch, inexplicably had comparable framerates on a high-end GPU, and these games were no doubt executing a great many non-linear instructions.

8 Conclusion

Implementing a basic computer (with an extant software library) using floating point addition and scaling demonstrates the highly general computing power they contain, despite approximating mathematically limited operations. We can say informally that they are Turing complete. This also renders the previous sections moot; performance notwithstanding, we could directly implement the Mandelbrot set, the \( \tanh \) transfer function, or AES using this 8-bit computer. It also immediately gives us a linear chess engine (including game tree search and a user interface) by emulating *chessmaster.nes*; in fact this engine already participated in our tournament (Figure 8)!

8.1 Future work

If I remember correctly (and I probably don’t), Gödel showed that an axiomatic system with addition and multiplication can encode sufficient facts about the natural numbers to engender *incompleteness* [9]. However, a system with only addition (such as Presburger arithmetic) does not have this problem. Incompleteness is similar to the halting problem for Turing complete systems, in that it is easy to encounter given a small set of primitives and the canonical demonstration is a diagonalization argument. Is floating point addition alone Turing complete? Can we prove it? If so, is the fact that real mathematical addition and multiplication have this deep incompleteness property related to the fact that IEEE-754 addition and multiplication have the deep computational property? Coincidence?

If not addition alone, the FMA (fused multiply-add) instruction very likely suffices, as it performs both a multiplication and addition. This makes sense, as the equation \( F = MA \) is fundamental to physics.

Linear logic???

Thinking about the 2A03 implementation, each loop executes the exact same set of instructions, with a high degree of parallelism. The use of condition codes mimics the way that VLIW machines and modern GPUs execute data-parallel programs. This seems to lend itself to highly parallel execution on GPUs; in fact the “Tensor cores”

---

21 No.

22 Yes.
designed for accelerating ML inference can likely execute these floating-point operations. Moreover, since the operations being executed are linear, the entire computation is trivially differentiable. This means that, if you don’t think about it too hard (but you need to think about it a medium amount of hard, because it is a confusing thought), you could use a finite sequence of NES instructions as transfer functions in a network, and back-propagate errors (giving an error vector towards a machine state and controller inputs that would yield the desired output state). This of course would not actually work, similar to how automatic differentiation does not actually work.

Not everyone uses IEEE-754 floating point these days. For example the bfloat16 format has gained traction in machine learning. Are similar tricks possible in these alternate universes, or is IEEE-754 simply the best forever?

Other applications of this technology are possible, and further study is warranted. For example, a common act in video editing is to rearrange clips from a source video in alphabetical order [2]. It was formerly believed that this required non-linear video editing (aside from “Already Being Filmed In Lexicographic Order Type Videos”). But it seems straightforward to use techniques from this paper to perform them linearly.

8.1.1 Conclusion Conclusion

A line has been drawn in the sand. \( y \) is truly equal to \( mx \) plus \( b \). The professor has been defeated. The dead horse has been beaten. The paper is finally over.

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