Introduction to Neural Networks
U. Minn. Psy 5038
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Lecture 4

Review: The generic connectionist model of the neuron

■ **Review**. In lecture 3, will developed some *Mathematica* tools to model Stage 1 and 2 of the generic connectionist neuron. Lists are particularly important. Vectors are lists of scalars that will be used to represent patterns of neural activity as well as synaptic weights of a neuron. Matrices are lists of vectors that will be used to represent a collection of synaptic weights over many model neurons.
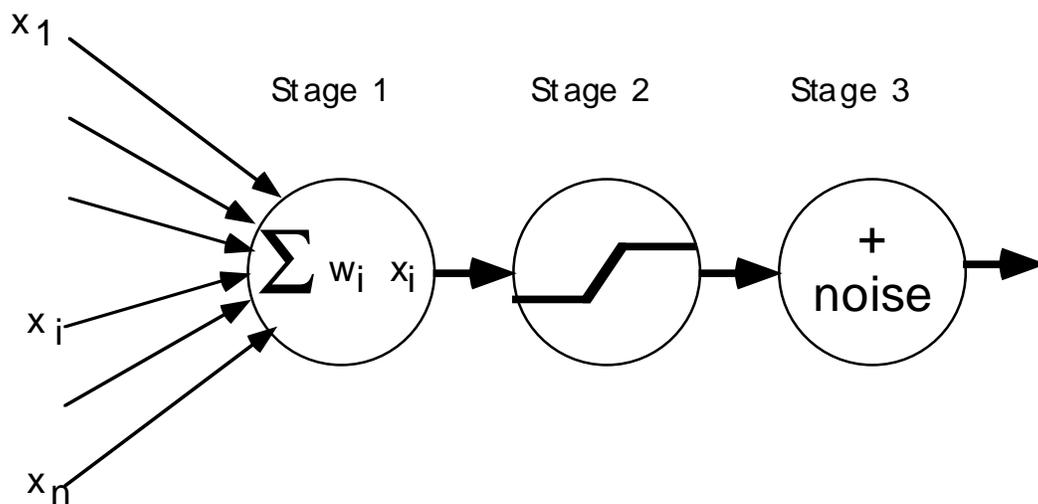
$$y = \sum_j w_j x_j$$

A single neuron

■ Stage 1: Vector dot product: Multiplying a vector of synaptic weights by a vector of input activities

In *Mathematica*, this linear weighted sum is written as a dot product:

```
w={w1,w2}; x = {x1,x2};
y = w.x
```

```
w1 x1 + w2 x2
```

■ Stage 2: The point non-linearity, squash functions.

$$y = \sigma \left\{ \sum_j w_j x_j \right\} + n$$

You've already seen how to define your own function for the sigmoidal non-linearity, $\sigma\{\}$.

```
squash[x_] := N[1/(1 + Exp[-x])];
```

Another form for a squashing function which is popular among physicists is:

```
squash2[x_] := ArcTan[x];
```

The values of **ArcTan[]** range from -Pi/2 to +Pi/2:

```
squash2[-Infinity]
```

$$-\frac{\pi}{2}$$

We've used *Mathematica* 's symbolic capabilities to confirm the lower range. Now run the output through the squash function:

```
y=squash[w.x];
```

Below we'll add stage 3, the noise term.

■ **Optional Exercise:** Plot squash[] and squash2[] on the same graph for values of x ranging from -5 to 5.
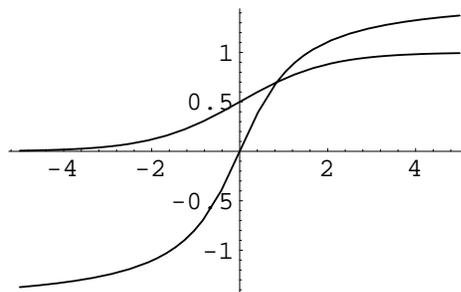
Either of the squashing functions can be used to model the small-signal compression and large signal saturation characteristics of neural output.

A two-layer neural network

Let's put the pieces together and model a simple two-layer neural network. We model the pattern of input activity by vector **x** with four input values, a 4x4 set of synaptic weights by matrix **W**, and the output pattern by vector **y**.

■ **Stage 1: Multiplying a matrix of synaptic weights by a vector of input activities**
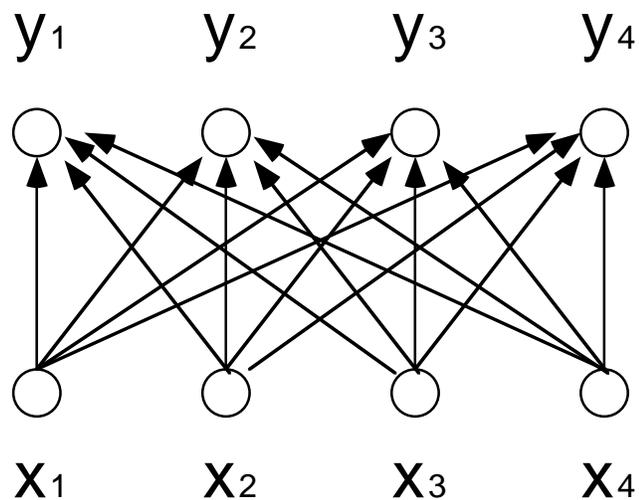
$$y_i = \sum_j w_{ij} x_j$$

$$y_i = \sigma \left\{ \sum_j w_{ij} x_j \right\}$$

In terms of *Mathematica*, let's assign positive random inputs and weights:

```
x = Table[Random[], {x,1,4}];
W = Table[Random[], {i,1,4},{j,1,4}];
y=W.x
```

{0.496035, 1.51142, 0.818805, 0.993001}

■ Stage 2: Point non-linearity

Recall in lecture 2, we pointed out that scalar functions are by default listable, which means that **squash[]** will get applied to each element of the vector **W.x** in turn. The output of a two-layer generic neural network can be written very concisely:

```
y = squash[W.x]
```

{0.621527, 0.819272, 0.693983, 0.72968}

■ Optional Exercise

Define **x** and **W** to model a two-layer neural network with 6 inputs and 2 outputs.

Modeling noise: Generic neuron plus noise

We'd like to add a Stage 3 to our model of the neuron in which we take account of the noisiness of neural transmission. For this, we need the notion of a *probability distribution*. We could develop the routines we need using basic *Mathematica* functions. However, much of the work has been done for us in the *Standard Mathematica Packages*(Add-ons, in the Help menu). These packages have to be read in when you need the function definitions they contain. As a first approximation the maintained action potential discharge can be modeled as a Poisson distribution. But to use the Poisson distribution in a *Mathematica* model, you have to read in the Statistics package **DiscreteDistributions** as shown below.

Statistics and stochastic processes

Statistical routines are useful for both theoretical aspects of modeling as well as for Monte Carlo simulations. So it is worth a little effort to get acquainted with some fundamental tools and definitions. Let's start by reading in one of the statistics packages and defining a Poisson distribution with a mean of 50 (e.g. 50 spikes per second of a neuron).

■ Discrete distributions

In[1]:= `<<Statistics`DiscreteDistributions``

In[2]:= `pdist = PoissonDistribution[50];`

The probability distribution function is given by:

In[3]:= `PDF[pdist,a]`

Out[3]= $\dfrac{50^a}{e^{50}\, a!}$

The output shows *Mathematica*'s definition of the function. You can obtain the mean, variance and standard deviation (which is the square root of the variance) of the distribution we've defined. Try it:

```
Mean[pdist]
Variance[pdist]
StandardDeviation[pdist]
```

What is your guess of the general relationship between the mean and variance for the Poisson distribution?

We are going to approximate the noisiness of neural discharge with a Normal or Gaussian distribution. The Gaussian distribution is continuous, rather than discrete. It is a fairly good approximation of a Poisson distribution for large values of the mean. To model the Gaussian, we need to read in the following package:

■ Continuous distributions, probability densities

In[4]:= `<<Statistics`ContinuousDistributions``

In[5]:= `ndist = NormalDistribution[0.5,.1];`

General::spell1 : Possible spelling error: new symbol name "ndist" is similar to existing symbol "pdist".

```
Print[Mean[ndist],", ",Variance[ndist],", ",
    StandardDeviation[ndist]]
```

0.5, 0.01, 0.1

A plot of the probability distribution function for this normal distribution looks like:

In[6]:= `Plot[PDF[ndist,x],{x,-.25,1.5}, PlotRange->{0,4},AxesLabel->{"x","p"}];`
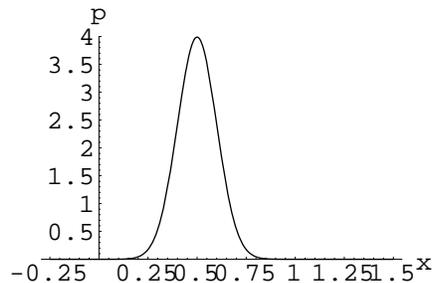
A given ordinate value is a "density", rather than probability. But we can talk about the probability that x takes on some value in an interval, say dx. For a small interval, dx, the probability $\simeq$ p(x)dx. What is the probability that x takes on some value between $+\infty$ and $-\infty$? What is area under this curve?

The cumulative distribution, f(x2) =p(x<x2), tells us the probability of x being less than a particular value of x2:

In[7]:= `Plot[CDF[ndist,x2],{x2,-.25,1.5},AxesLabel->{"x2","f"}];`
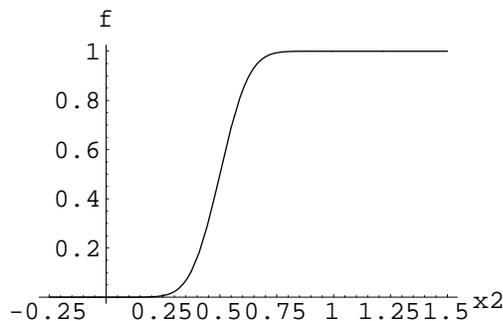
You can see from the graph that for this distribution, once x2 is greater than 0.7 or so, the probability of x being less than that is virtually certain, i.e. is essentially 1.

■ Statistical Sampling

Having defined the normal distribution, how can we draw samples from it? In other words, can we simulate a process in which we fill a hat with slips of paper in such a way that the proportions for each value mimic what we obtain from a theoretical distribution?

Most standard programming languages come with subroutines for doing pseudo-random number generation. Unlike the Poisson or Gaussian distribution, these numbers are usually **uniformly distributed--**that is, the probability of being a certain value (or within a tiny range) is constant over the entire sampling range of the random variable.

This is like filling the hat with slips of paper where the number of slips is the same for each value.

Mathematica comes with a standard function, **Random[]** that enables us to generate random numbers that are uniform, Poisson, Normal, depending on the argument. (There are some other possible distributions in the packages too, like the **ChiSquareDistribution**).

```
??Random
```

Random[ ] gives a uniformly distributed pseudorandom Real in the range 0 to 1. Random[
     type, range] gives a pseudorandom number of the specified type, lying in the specified
     range. Possible types are: Integer, Real and Complex. The default range is 0 to 1. You can
     give the range {min, max} explicitly; a range specification of max is equivalent to {0, max}.
     Random[distribution] gives a random number with the specified statistical distribution.

Attributes[Random] = {Protected}

```
Random[ndist]
```

0.701706

Putting together stages 1, 2 and 3 together

We can do everything at once, producing the output of a generic neuron, with synaptic weights **w**, neural noise with a mean of 0.0 and std. dev. 0.1 to an input x:

```
w = {2,1,-2,3};
ndist2 = NormalDistribution[0.0,.1];
y[x1_] := N[squash[w.x1] + Random[ndist2]];
y[{2,3,0,1}]
```

0.911057

If we invoke the **y[]** function again, we get a different response:

```
y[{2,3,0,1}]
```

0.890749

To sum up, the model you should have in mind is that at any given time interval (which is implicit in this continuous-response, discrete-time model), the neuron computes the sum of its weighted inputs, and the output signal, y, is a spike rate over this interval. With a sigmoidal non-linearity, there is small-signal suppression, and large-signal saturation.

■ Exercise

Suppose all the inputs except the first are clamped at zero. What does the response, **y[{t,0,0,0}]**, look like as a function of x for various levels of noise?

```
Plot[y[{x, 0, 0, 0}], {x, -2, 2}, PlotRange -> {0, 2},
 AxesLabel -> {"Input signal, x", "Frequency"}];
```
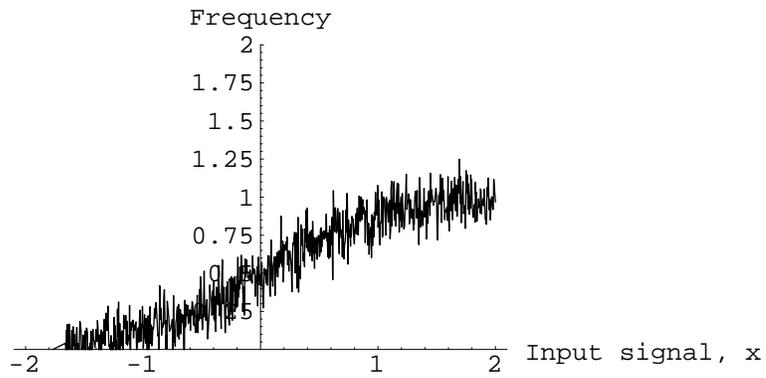
Vector operations and patterns of neural activity

■ **State space and state vectors.**

■ In neural networks, we are often concerned with a vector whose components represent the activities of neurons which are changing in time. So sometimes we will talk about **state vectors**.There isn't anything profound about this terminology--it just reflects that we are interested in the value of the vector when the system is in a particular state at time t. It is often very useful to think of an n-dimensional vector as a point in an n-dimensional space. This space is often referred to as **state space**. Suppose, we have a 3 neuron system. We can describe the state of this system as a 3-dimensional vector where each component represents the activity of the neuron. Further, suppose just for the sake of an example to visualize, the activities of the first, second, and third neurons (i.e. components) of a 3-dimensional vector are given by: **y = {Cos[t], Sin[t], t}**. We can use the Mathematica function, **ParametricPlot3D[]** to get a picture of how this state vector evolves in time through state space:

```
Clear[y];
y[t_]  :=  {Cos[t], Sin[t], Cos[2 t]};
ParametricPlot3D[y[t], {t,0,5}, AxesLabel->{"x","y","z"}];
```

ParametricPlot3D::ppcom :
    Function *y*(*t*) cannot be compiled; plotting will proceed with the uncompiled function.
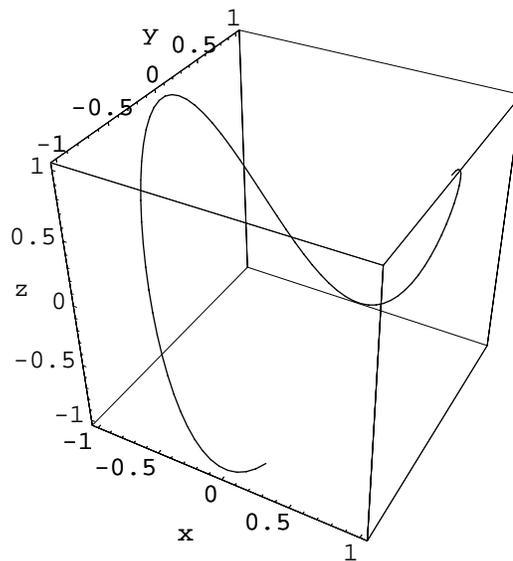
■ **Dimension of a vector.**

You can get the dimensionality of a vector using **Dimensions[]**, or **Length[]**.

In[8]:=
```
v = {2.1, 3, -0.45, 4.9};

Dimensions[v]
```

Out[9]=
```
{4}
```

Dimensions[], will give you the dimensions of a matrix, while **Length[]** tells you the number of elements in the list. For example,

```
M = {{2,4,2}, {1,6,4}};
```

```
Length[M]
```

```
2
```

Try comparing **Length[M]** with **Dimensions[M]**.

■ **Transpose of a vector.**

The transpose of a column vector is just the same vector arranged in a row. However, because of the way Mathematica uses lists to represent vectors you don't have to distinguish between row and column vectors. In standard math notation, transpose of a vector **x**, is often written $\mathbf{x}^T$. You can see a vector in column form by typing **v//MatrixForm**, or**:**

In[13]:=
```
MatrixForm[v]
```

Out[13]//MatrixForm=

$$\begin{pmatrix} 2.1 \\ 3 \\ -0.45 \\ 4.9 \end{pmatrix}$$

■ **Vector addition** is accomplished by simply adding the components of each vector to make a new vector. Note that the vectors all have the same dimension.

```
a = {3,1,2};
b = {2,4,8};
c = a + b
```

```
{5, 5, 10}
```

Vectors can be multiplied by a constant. We saw an example of this earlier.

```
2 a
```

{6, 2, 4}

■ Metric length of a vector.

It is unfortunate terminology, but **Length[]** does NOT give you the metrical length of the vector. In order to get the length of a vector, you calculate the Euclidean distance from the origin to the end-point of the vector. We get this by squaring each component, adding up the squares, and taking the square root. First, we will do this using the **Apply[]** function, where the **Plus** operation is applied to all the elements of the list. Note that the operation of exponentiation is "listable", that is it is applied to each element of the vector:

```
a^2
```

{9, 1, 4}

What is **a a** ?

```
N[Sqrt[Apply[Plus, a^2]]]
```

3.74166

If you wish, you can define your own function to apply to the list. What we have just calculated is the square root of the dot product or inner product of **a** with itself. The length of a vector **a** is often written as |**a**| in standard math notation. In the next section, we use the inner or dot product to calculate the metric length of a vector.

■ Inner product. To calculate the inner product of two vectors, you multiply the corresponding components and add them up:

```
u = {u1,u2,u3,u4};
v = {v1,v2,v3,v4};
u.v
```

u1 v1 + u2 v2 + u3 v3 + u4 v4

The **inner product** is also called the **dot product**. Later we will see what is meant by **outer product**. The inner product between two vectors **a** and **b** is written either as:

$$\mathbf{a.b} \text{ or } [\mathbf{a,b}], \text{ or } \mathbf{a^T b}$$

Mathematica uses the dot notation.

One use of the inner product is to calculate the length of a vector. **a.a** is just the sum of the squares of the elements of **a**, so gives us another way of calculating the length of a vector.

```
N[Sqrt[a.a]]
```

```
 3.74166
```

Let's define a function that will return the length of a vector, x:

```
Vectorlength[x_]  :=  N[Sqrt[x.x]]
```

■ **Projection.** The dot product, **a.b**, is equal to:

$$|a|\ |b|\ cos(\text{angle between } a \text{ and } b)$$

In problem set 1, you calculate the output of a linear neuron model as the dot product between an input vector and a weight vector. Both the weight and input lists can be thought of as vectors in an n-dimensional space. Suppose the weight vector has unit length. Recall that you can normalize any vector to unit length by dividing by its length:

```
v = v/Sqrt[v.v] ;
```

Geometrically, we can think of the output of a neuron as the projection of the activity of the neuron input activity vector onto the weight vector direction. Suppose the input vector is already perpendicular to the weight vector, then the output of the neuron is zero, because the cosine of 90 degrees is zero. As you found or will find with the cross-correlator of Problem Set 1, the further the input pattern is away from the weight vector, as measured by the cosine between them, the poorer the "match" between input and weight vectors, and the lower the response.

Here are three lines of code that calculate the two-dimensional vector **z** in the direction of **w**, with a length determined by "how much of **x** projects in the **w** direction":

```
x = {1,2};
w = N[{2/Sqrt[5],1/Sqrt[5]}];
z = (x.w) w ;
```

```
Show[   Graphics[{Line[{{0,0}, x}],
            Line[{{0,0}, z}],
            {Dashing[{0.03,0.03}], Line[{x, z}] },
            Text[FontForm["w",{"Helvetica-Bold",18}], w, {0,-1}],
            Text[FontForm["x",{"Helvetica-Bold",18}], x, {-2,0}],
            Text[FontForm["z",{"Helvetica-Bold",18}], z, {0,-1}],
            {AbsoluteThickness[3], Line[{{0,0}, w}] }
        }],
        Axes->True, AspectRatio->1

];
```
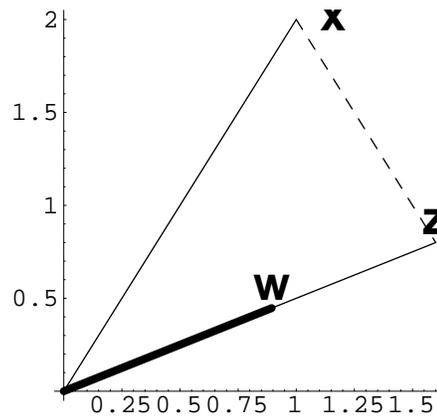

■ Angle between two vectors and orthogonality: Similarity measure between patterns

Often we will want some measure of the similarity between two patterns of neural firings. As we have just seen, one measure of comparison is the degree to which the two state vectors point in the same direction. The cosine of the angle between two vectors is one possible measure:

```
cosine[x_,y_] := x.y/(Vectorlength[x] Vectorlength[y])
cosine[a,b]
```

> 0.758175

Note that if two vectors point in the same direction, the cosine of the angle between them is 1:

```
a = {2,1,3,6};
b = {6, 3, 9, 18};
cosine[a,b]
```

> 1.

Try verifying that **w** and **z** from the previous section point in the same direction.

If two vectors point in the opposite directions, the cosine of the angle between them is -1:

```
a = {-2,-1,-3,-6};
b = {6, 3, 9, 18};
cosine[a,b]
```

$-1.$

Two vectors may point in the same direction, but could be quite different because they have different lengths. Another measure of similarity is the length of the difference between two vectors:

```
Vectorlength[a - b]
```

28.2843

■ **Orthogonality.** The case where vectors are at right angles to each other is an important special case that is worth spending a little time on. Consider an 8-dimensional space. One very familiar set of orthogonal vectors is the following:

```
u1 = {1,0,0,0,0,0,0,0};
u2 = {0,1,0,0,0,0,0,0};
u3 = {0,0,1,0,0,0,0,0};
u4 = {0,0,0,1,0,0,0,0};
u5 = {0,0,0,0,1,0,0,0};
u6 = {0,0,0,0,0,1,0,0};
u7 = {0,0,0,0,0,0,1,0};
u8 = {0,0,0,0,0,0,0,1};
```

Each vector has unit length, and it is easy to see just by inspection that the inner product between any two is zero. On the other hand, here is another set of 8 vectors in 8-space for which it is not immediately obvious that they are all orthogonal. These vectors are called Walsh functions:

```
v1 = {1, 1, 1, 1, 1, 1, 1, 1};
v2 = {1,-1,-1, 1, 1,-1,-1, 1};
v3 = {1, 1,-1,-1,-1,-1, 1, 1};
v4 = {1,-1, 1,-1,-1, 1,-1, 1};
v5 = {1, 1, 1, 1,-1,-1,-1,-1};
v6 = {1,-1,-1, 1,-1, 1, 1,-1};
v7 = {1, 1,-1,-1, 1, 1,-1,-1};
v8 = {1,-1, 1,-1, 1,-1, 1,-1};
```

You can calculate the inner products between any two, and you will find out that they are all zero. Note that with the first set of vectors, $\{\mathbf{u_i}\}$, you can tell which vector it is just by looking for where the 1 is. For the second set, $\{\mathbf{v_i}\}$, you can't tell by looking at just one component. For example, the first component of all of the Walsh functions has a 1. You have to look at the pattern to tell which Walsh function you are looking at.

Suppose for the moment that we want to assign meaning to each of the patterns--each pattern is a code for some thing, like "grandma Tompkins", "grandma Wilke", and so forth. If we use the **u**'s, then we could look for the one neuron that lights up

to find out which grandma it is representing--then neuron activity represented, for example, by the third element of the pattern could mean "grandma Wilke". This strategy wouldn't work if we encoded a collection of grandmothers using the **v**'s. The **v**'s give us a simple example of what is sometimes referred to as a **distributed code**. The **w**'s are examples of a **grandmother cell code**. The reason for this obscure terminology can be traced to earlier debates on whether there may be single cells in the brain whose firing uniquely determines the recognition of one's grandmother.

■ **Orthonormality.** The Walsh set is orthogonal, but they are not of unit length. We have already seen some of the advantages of working with unit length vectors. The general issue of normalization comes up all the time in neural networks both in terms of limiting overall neural activity, and limiting synaptic weights. So it is sometimes convenient to normalize an orthogonal set, producing what is known as an *orthonormal* set of vectors:

```
w1 = v1/Vectorlength[v1];
w2 = v2/Vectorlength[v2];
w3 = v3/Vectorlength[v3];
w4 = v4/Vectorlength[v4];
w5 = v5/Vectorlength[v5];
w6 = v6/Vectorlength[v6];
w7 = v7/Vectorlength[v7];
w8 = v8/Vectorlength[v8];
```

Vector representations, linear algebra

The issue of how information is to be represented is fundamental in the information sciences generally, as well as for neural network theory. A pattern of activity over a set of neurons is presumed to mean something, and there are different ways of coding the same meaning. But different codes have different properties. A code may not be sufficient to uniquely code all the possible things we need to represent. A code could be redundant and have more than one way of representing the same thing. This section continues with our review of the basics of vector and linear algebra by going a little more deeply into the subject. The pay-off will be some mathematics that provides intuition about issues of neural representation. You can think of this as a first lesson in the "psychology of linear algebra".

■ **Basis sets**

It is pretty clear that given any vector whatsoever in 8-space, you can specify how much of it gets projected in each of the eight directions specified by the unit vectors v1, v2, ...v8. But you can also build back up an arbitrary vector by adding up all the contributions from each of the component vectors. This is a consequence of vector addition and can be easily seen to be true in 2 dimensions. We can verify it ourselves. Pick an arbitrary vector g, project it onto each of the basis vectors, and then add them back up again:

```
g = {2,6,1,7,11,4,13, 29} ;
```

```
(g.u1) u1   + (g.u2) u2  +(g.u3) u3  +(g.u4) u4   +
(g.u5) u5   +(g.u6) u6  +(g.u7) u7  +(g.u8) u8
```

```
{2, 6, 1, 7, 11, 4, 13, 29}
```

■ Exercise

What happens if you project **g** onto the normalized Walsh basis set defined by **{w1,w2,...}** above, and then add up all 8 components?

```
(g.w1) w1   + (g.w2) w2  +(g.w3) w3  +(g.w4) w4   +
(g.w5) w5   +(g.w6) w6  +(g.w7) w7  +(g.w8) w8
```

The projections, **g.u$_i$** are sometimes called the **spectrum** of **g**. This terminology comes from the Fourier basis set used in Fourier analysis. A discrete version of a Fourier basis set is similar to the Walsh set, except that the elements fit a sine wave pattern, and so are not binary-valued.

The orthonormal set of vectors we've defined above is said to be **complete**, because any vector in 8-space can be expressed as a linear weighted sum of these **basis vectors**. The weights are just the projections. If we had only 7 vectors in our set, then we would not be able to express any 8-dimensional vector in terms of this basis set. The seven vector set would be said to be **incomplete**. A basis set which is orthonormal and complete is very nice from a mathematical point of view. Another bit of terminology is that these seven vectors would not **span** the 8-dimensional space. But they would span some sub-space, that is of smaller dimension, of the 8-space.

There has been much interest in describing the effective weighting properties of visual neurons in primary visual cortex of higher level mammals (cats, monkeys) in terms of basis vectors. One issue is if the input (e.g. an image) is projected (via a collection of receptive fields) onto a set of neurons, is information lost? If the set of weights representing the receptive fields of the collection of neurons is complete, then no information is lost.

■ **Linear dependence**

What if we had 9 vectors in our basis set used to represent vectors in 8-space? For the u's, it is easy to see that in a sense we have too many, because we could express the 9th in terms of a sum of the others. This set of nine vectors would be said to be linearly dependent. A set of vectors is linearly dependent if one or more of them can be expressed as a linear combination of some of the others. Sometimes there is an advantage to having an "over-complete" basis set (e.g. more than 8 vectors for 8-space; cf. Simoncelli et al., 1992).

Theorem: A set of mutually orthogonal vectors is linearly independent.

However, note it is quite possible to have a linearly independent set of vectors which are not orthogonal to each other. Imagine 3-space and 3 vectors which do not jointly lie on a plane. This set is linearly independent.

If we have a linearly independent set, say of 8 vectors for our 8-space, then no member can be dropped without a loss in the dimensionality of the space spanned.

It is useful to think about the meaning of linear independence in terms of geometry. A set of three linearly independent vectors can completely span 3-space. So any vector in 3-space can be represented as a weighted sum of these 3. If one of the members in our set of three can be expressed in terms of the other two, the set is not linearly independent and the set only spans a 2-dimensional subspace. That is, the set can only represent vectors which lay on a plane in 3-space. This can be easily seen to be true for the set of u's, but is also true for the set of v's.

■ **Thought exercise**

Suppose there are three inputs feeding into three neurons in the simple linear network such as defined at the beginning of this lecture. If the weight vectors of the three neurons are not linearly independent, do we lose information?

Linearity, real neural networks, and what's up next time?

From a computational standpoint, the squashing function has both advantages and disadvantages. It is what makes our neural network model *non-linear*, and as we will see later, this non-linearity enables networks to compute functions that can't be computed with a linear network. On the other hand, non-linearities make the analysis complicated. In fact, there are cases for which most of the neural activities are in the mid-range of the squashing function, and here one can approximate the network as a purely linear one--just matrix operations on vector inputs, and the analysis becomes relatively simple.

Compared to the complexity of real neurons and networks, assuming linearity might seem to be just too simple. But we will see in the next lecture, that a linear model can be quite good model for some biological subsystems. We will apply the techniques of linear vector algebra to model a network discovered in the visual system of the horseshoe crab.

References

Simoncelli, E. P., Freeman, W. T., Adelson, E. H., & Heeger, D. J. (1992). Shiftable Multi-scale Transforms. <u>IEEE Trans. Information Theory</u>, <u>38</u>(2), 587--607.

Strang, G. (1988). <u>Linear Algebra and Its Applications</u> (3rd ed.). Saunders College Publishing Harcourt Brace Jovanovich College Publishers.