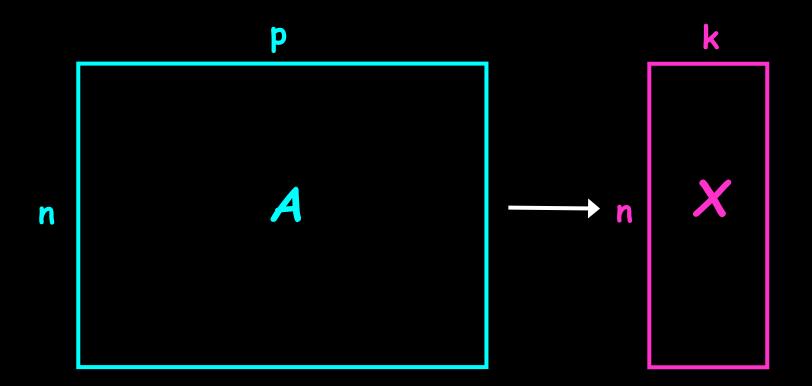
Principal Component Analysis (PCA)

Data Reduction

summarization of data with many (p)
variables by a smaller set of (k) derived
(synthetic, composite) variables.



Data Reduction

- "Residual" variation is information in A that is not retained in X
- balancing act between
 - clarity of representation, ease of understanding
 - oversimplification: loss of important or relevant information.

Principal Component Analysis (PCA)

- probably the most widely-used and wellknown of the "standard" multivariate methods
- invented by Pearson (1901) and Hotelling (1933)
- first applied in ecology by Goodall (1954) under the name "factor analysis" ("principal factor analysis" is a synonym of PCA).

Principal Component Analysis (PCA)

- takes a data matrix of n objects by p variables, which may be correlated, and summarizes it by uncorrelated axes (principal components or principal axes) that are linear combinations of the original p variables
- the first k components display as much as possible of the variation among objects.

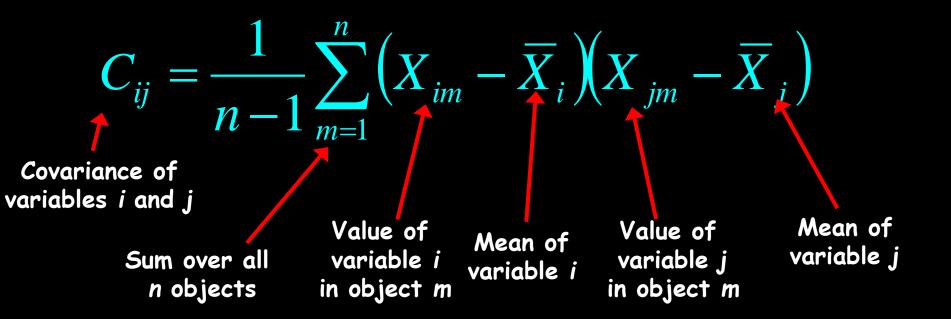
Geometric Rationale of PCA

- objects are represented as a cloud of n points in a multidimensional space with an axis for each of the p variables
- the centroid of the points is defined by the mean of each variable
- the variance of each variable is the average squared deviation of its n values around the mean of that variable.

$$V_i = \frac{1}{n-1} \sum_{m=1}^{n} (X_{im} - \overline{X}_i)^2$$

Geometric Rationale of PCA

 degree to which the variables are linearly correlated is represented by their covariances.

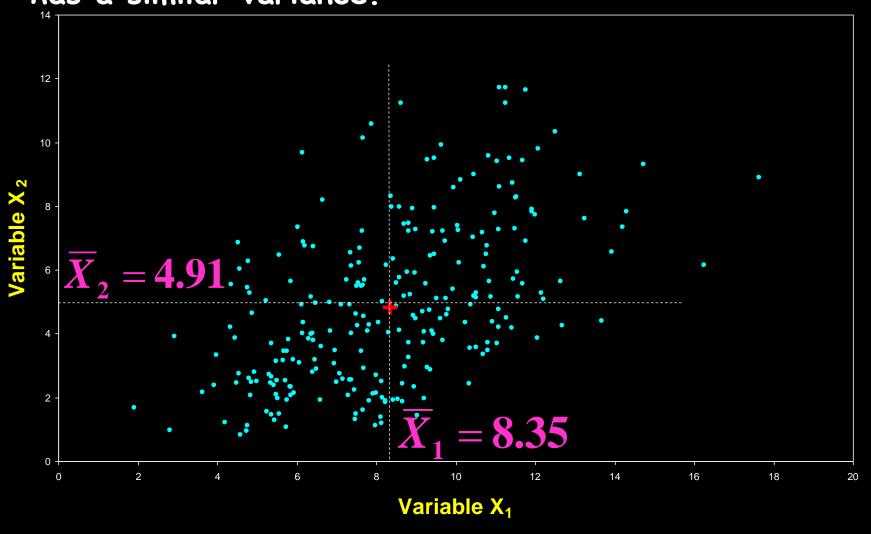


Geometric Rationale of PCA

- objective of PCA is to rigidly rotate the axes of this p-dimensional space to new positions (principal axes) that have the following properties:
 - ordered such that principal axis 1 has the highest variance, axis 2 has the next highest variance,, and axis p has the lowest variance
 - covariance among each pair of the principal axes is zero (the principal axes are uncorrelated).

2D Example of PCA

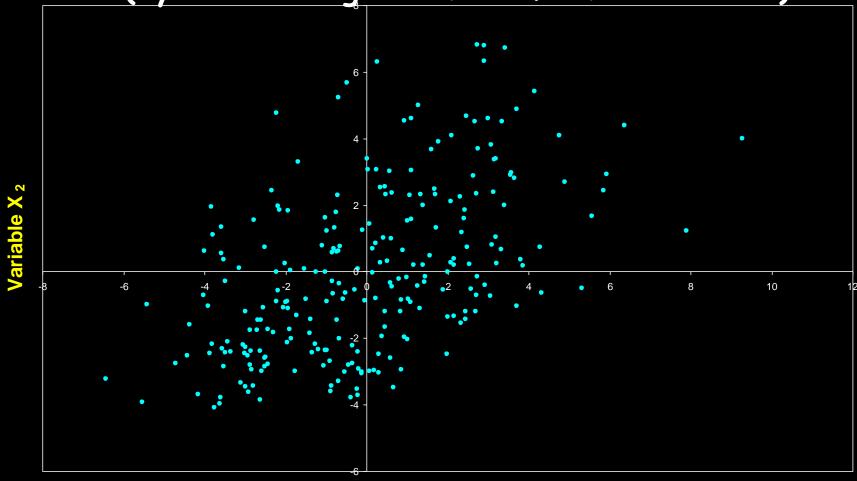
variables X_1 and X_2 have positive covariance & each has a similar variance.



 $V_1 = 6.67$ $V_2 = 6.24$ $C_{1,2} = 3.42$

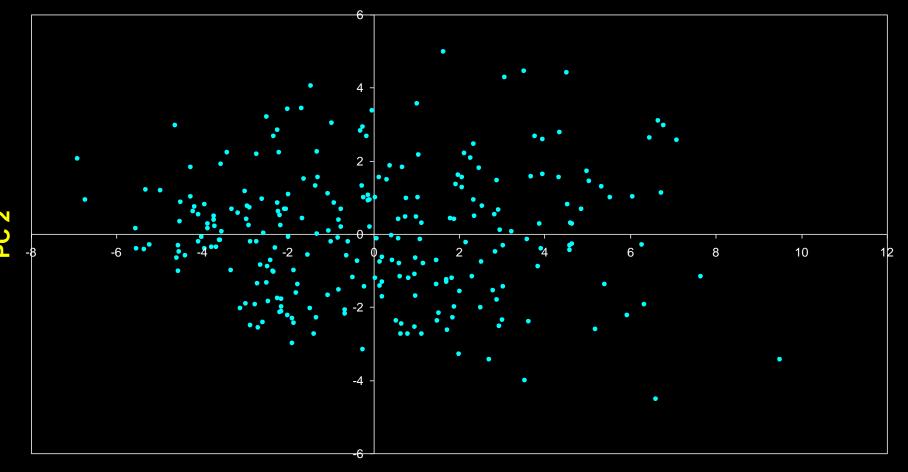
Configuration is Centered

• each variable is adjusted to a mean of zero (by subtracting the mean from each value).



Principal Components are Computed

- PC 1 has the highest possible variance (9.88)
- PC 2 has a variance of 3.03
- PC 1 and PC 2 have zero covariance.



The Dissimilarity Measure Used in PCA is Euclidean Distance

- PCA uses Euclidean Distance calculated from the p variables as the measure of dissimilarity among the n objects
- PCA derives the best possible k dimensional (k < p) representation of the Euclidean distances among objects.

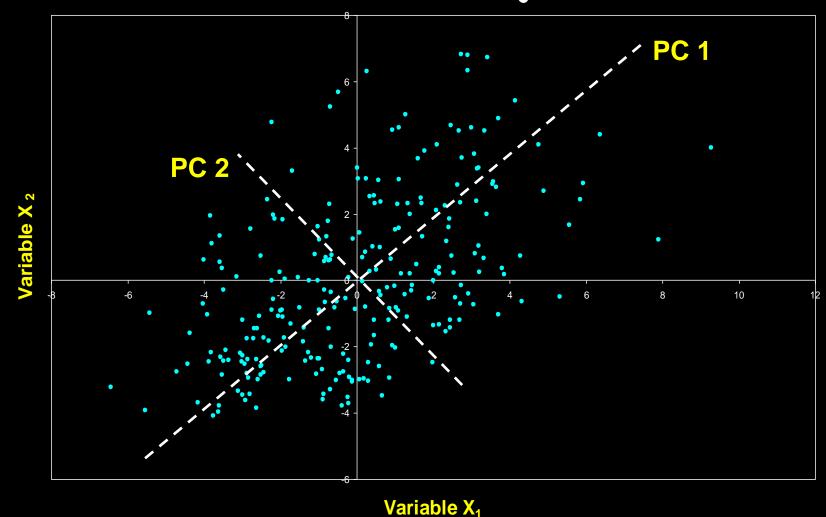
Generalization to p-dimensions

- In practice nobody uses PCA with only 2 variables
- The algebra for finding principal axes readily generalizes to p variables
- PC 1 is the direction of maximum variance in the p-dimensional cloud of points
- PC 2 is in the direction of the next highest variance, subject to the constraint that it has zero covariance with PC 1.

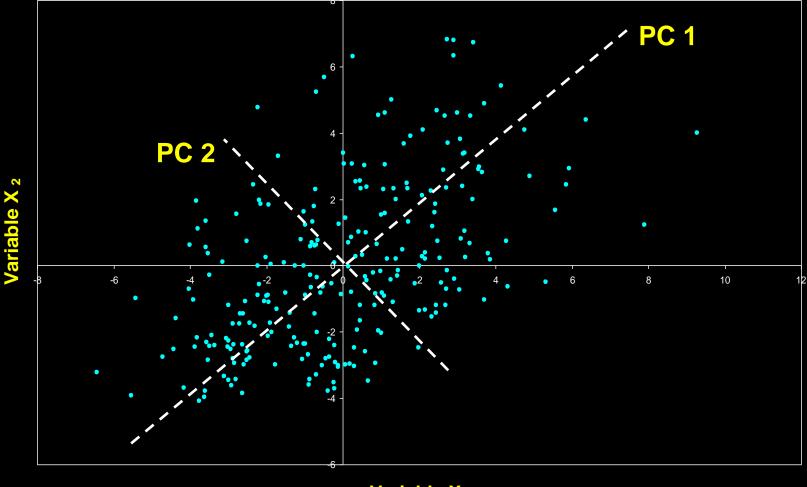
Generalization to p-dimensions

- PC 3 is in the direction of the next highest variance, subject to the constraint that it has zero covariance with both PC 1 and PC 2
- and so on... up to PC p

- each principal axis is a linear combination of the original two variables
- $PC_j = a_{i1}Y_1 + a_{i2}Y_2 + ... a_{in}Y_n$ a_{ij} 's are the coefficients for factor i, multiplied by the measured value for variable j



- PC axes are a rigid rotation of the original variables
- PC 1 is simultaneously the direction of maximum variance and a least-squares "line of best fit" (squared distances of points away from PC 1 are minimized).



Generalization to p-dimensions

- if we take the first k principal components, they define the k-dimensional "hyperplane of best fit" to the point cloud
- of the total variance of all p variables:
 - PCs 1 to k represent the maximum possible proportion of that variance that can be displayed in k dimensions
 - i.e. the squared Euclidean distances among points calculated from their coordinates on PCs 1 to k are the best possible representation of their squared Euclidean distances in the full p dimensions.

Covariance vs Correlation

- using covariances among variables only makes sense if they are measured in the same units
- even then, variables with high variances will dominate the principal components
- these problems are generally avoided by standardizing each variable to unit variance and zero mean.

$$m{X}_{im}' = rac{\left(m{X}_{im} - m{\overline{X}}_i
ight)^{ ext{Mean}}}{ ext{SD}_i}$$
 Standard deviation of variable i

Covariance vs Correlation

- covariances between the standardized variables are correlations
- after standardization, each variable has a variance of 1.000
- correlations can be also calculated from the variances and covariances:

```
Correlation between variables i and j
r_{ij} = \frac{C_{ij}}{V_i V_j}
Variance of variable i
Variance of variable j
of variable i
```

- The Algebra of PCA

 first step is to calculate the crossproducts matrix of variances and covariances (or correlations) among every pair of the p variables
- square, symmetric matrix
- · diagonals are the variances, off-diagonals are the covariances.

	X ₁	X ₂
X ₁	6.6707	3.4170
X ₂	3.4170	6.2384

	X ₁	X ₂
X ₁	1.0000	0.5297
X ₂	0.5297	1.0000

Variance-covariance Matrix

Correlation Matrix

· in matrix notation, this is computed as

$$S = X'X$$

 where X is the n x p data matrix, with each variable centered (also standardized by SD if using correlations).

	X ₁	X ₂
X ₁	6.6707	3.4170
X ₂	3.4170	6.2384

	X ₁	X ₂
X ₁	1.0000	0.5297
X ₂	0.5297	1.0000

Variance-covariance Matrix

Correlation Matrix

Manipulating Matrices

 transposing: could change the columns to rows or the rows to columns

$$X = \begin{bmatrix} 10 & 0 & 4 \\ 7 & 1 & 2 \end{bmatrix}$$
 $X' = \begin{bmatrix} 10 & 7 \\ 0 & 1 \\ 4 & 2 \end{bmatrix}$

- multiplying matrices
 - must have the same number of columns in the premultiplicand matrix as the number of rows in the postmultiplicand matrix

- sum of the diagonals of the variancecovariance matrix is called the trace
- it represents the total variance in the data
- it is the mean squared Euclidean distance between each object and the centroid in p-dimensional space.

	X ₁	X ₂
X ₁	6.6707	3.4170
X ₂	3.4170	6.2384

	X ₁	X ₂
X ₁	1.0000	0.5297
X ₂	0.5297	1.0000

Trace = 12.9091

Trace = 2.0000

- finding the principal axes involves eigenanalysis of the cross-products matrix (5)
- the eigenvalues (latent roots) of S are solutions (λ) to the characteristic equation

$$|S - \lambda I| = 0$$

- the eigenvalues, λ_1 , λ_2 , ... λ_p are the variances of the coordinates on each principal component axis
- the sum of all p eigenvalues equals the trace of S (the sum of the variances of the original variables).

	X ₁	X ₂
X ₁	6.6707	3.4170
X ₂	3.4170	6.2384

$$\lambda_1 = 9.8783$$
 $\lambda_2 = 3.0308$

Note: $\lambda_1 + \lambda_2 = 12.9091$

- each eigenvector consists of p values which represent the "contribution" of each variable to the principal component axis
- eigenvectors are uncorrelated (orthogonal)
 - their cross-products are zero.

Eigenvectors			
	u ₁	u ₂	
X ₁	0.7291	-0.6844	
X ₂	0.6844	0.7291	

0.7291*(-0.6844) + 0.6844*0.7291 = 0

• coordinates of each object i on the k^{th} principal axis, known as the scores on PC k, are computed as

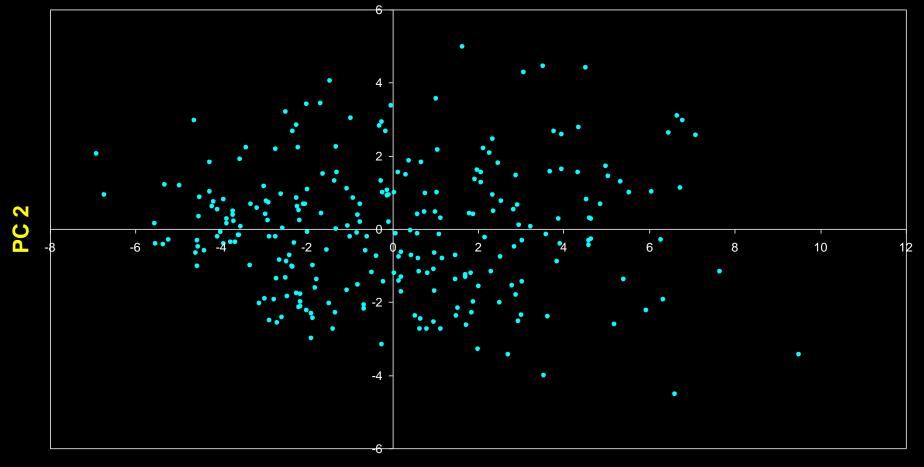
$$z_{ki} = u_{1k} x_{1i} + u_{2k} x_{2i} + \cdots + u_{pk} x_{pi}$$

• where Z is the n x k matrix of PC scores, X is the n x p centered data matrix and U is the p x k matrix of eigenvectors.

- variance of the scores on each PC axis is equal to the corresponding eigenvalue for that axis
- the eigenvalue represents the variance displayed ("explained" or "extracted") by the k^{th} axis
- the sum of the first k eigenvalues is the variance explained by the k-dimensional ordination.

 $\lambda_1 = 9.8783$ $\lambda_2 = 3.0308$ Trace = 12.9091

PC 1 displays ("explains")
9.8783/12.9091 = 76.5% of the total variance



- The cross-products matrix computed among the p principal axes has a simple form:
 - all off-diagonal values are zero (the principal axes are uncorrelated)
 - the diagonal values are the eigenvalues.

	PC ₁	PC ₂
PC ₁	9.8783	0.0000
PC ₂	0.0000	3.0308

Variance-covariance Matrix of the PC axes

A more challenging example

- data from research on habitat definition in the endangered Baw Baw frog
- 16 environmental and structural variables measured at each of 124 sites
- correlation matrix used because variables have different units



Eigenvalues

Axis	Eigenvalue	% of Variance	Cumulative % of Variance
1	5.855	36.60	36.60
2	3.420	21.38	57.97
3	1.122	7.01	64.98
4	1.116	6.97	71.95
5	0.982	6.14	78.09
6	0.725	4.53	82.62
7	0.563	3.52	86.14
8	0.529	3.31	89.45
9	0.476	2.98	92.42
10	0.375	2.35	94.77

Interpreting Eigenvectors

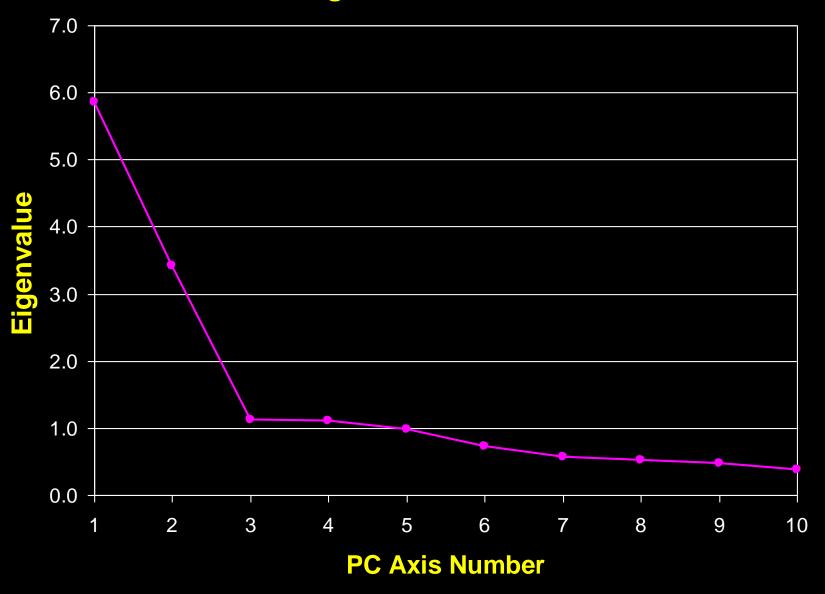
- correlations
 between variables
 and the principal
 axes are known
 as loadings
- each element of the eigenvectors represents the contribution of a given variable to a component

1	2	3
0.3842	0.0659	-0.1177
-0.1159	0.1696	-0.5578
-0.2729	-0.1200	0.3636
0.0538	-0.2800	0.2621
-0.0765	0.3855	-0.1462
0.0248	0.4879	0.2426
0.0599	0.4568	0.2497
0.0789	0.4223	0.2278
0.3305	-0.2087	-0.0276
-0.3053	0.1226	0.1145
-0.3144	0.0402	-0.1067
-0.0886	-0.0654	-0.1171
0.1364	-0.1262	0.4761
-0.3787	0.0101	0.0042
-0.3494	-0.1283	0.1166
0.3899	0.0586	-0.0175
	0.3842 -0.1159 -0.2729 0.0538 -0.0765 0.0248 0.0599 0.0789 0.3305 -0.3053 -0.3144 -0.0886 0.1364 -0.3787 -0.3494	0.3842 0.0659 -0.1159 0.1696 -0.2729 -0.1200 0.0538 -0.2800 -0.0765 0.3855 0.0248 0.4879 0.0599 0.4568 0.0789 0.4223 0.3305 -0.2087 -0.3053 0.1226 -0.3144 0.0402 -0.0886 -0.0654 0.1364 -0.1262 -0.3787 0.0101 -0.3494 -0.1283

How many axes are needed?

- does the $(k+1)^{th}$ principal axis represent more variance than would be expected by chance?
- several tests and rules have been proposed
- a common "rule of thumb" when PCA is based on correlations is that axes with eigenvalues > 1 are worth interpreting

Baw Baw Frog - PCA of 16 Habitat Variables

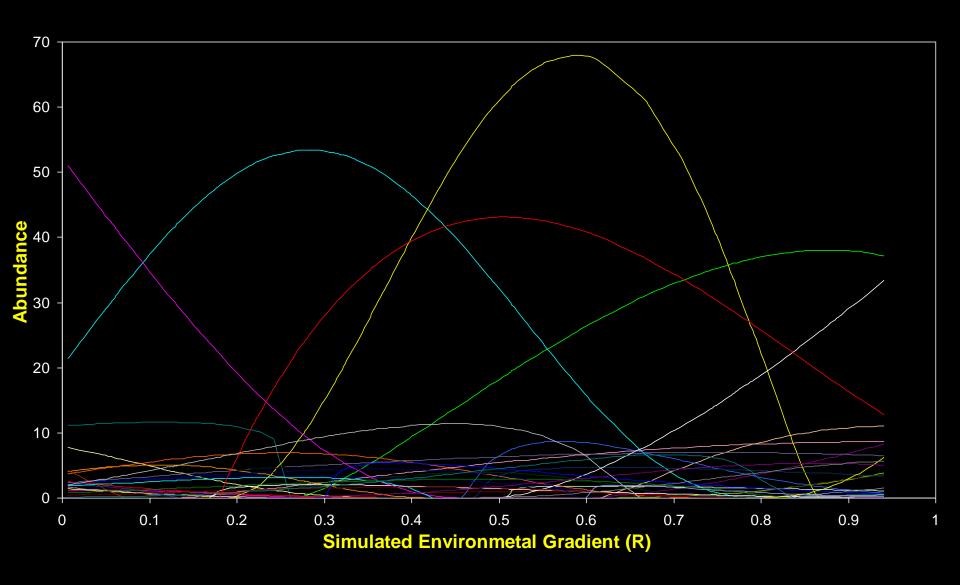


What are the assumptions of PCA?

- assumes relationships among variables are LINEAR
 - cloud of points in p-dimensional space has linear dimensions that can be effectively summarized by the principal axes
- if the structure in the data is NONLINEAR (the cloud of points twists and curves its way through p-dimensional space), the principal axes will not be an efficient and informative summary of the data.

When should PCA be used?

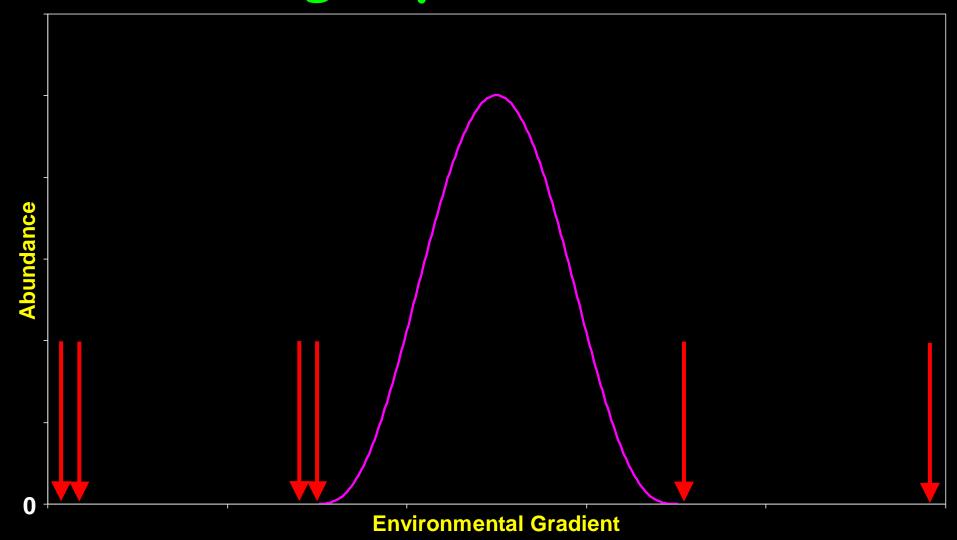
- In community ecology, PCA is useful for summarizing variables whose relationships are approximately linear or at least monotonic
 - e.g. A PCA of many soil properties might be used to extract a few components that summarize main dimensions of soil variation
- PCA is generally NOT useful for ordinating community data
- Why? Because relationships among species are highly nonlinear.



The "Horseshoe" or Arch Effect

- community trends along environmenal gradients appear as "horseshoes" in PCA ordinations
- none of the PC axes effectively summarizes the trend in species composition along the gradient
- SUs at opposite extremes of the gradient appear relatively close together.

Ambiguity of Absence



Beta Diversity 2R - Covariance Axis 2 Axis 1

The "Horseshoe" Effect

- curvature of the gradient and the degree of infolding of the extremes increase with beta diversity
- PCA ordinations are not useful summaries of community data except when beta diversity is very low
- using correlation generally does better than covariance
 - this is because standardization by species improves the correlation between Euclidean distance and environmental distance.

What if there's more than one underlying ecological gradient?

The "Horseshoe" Effect

- when two or more underlying gradients with high beta diversity a "horseshoe" is usually not detectable
- the SUs fall on a curved hypersurface that twists and turns through the p-dimensional species space
- · interpretation problems are more severe
- PCA should NOT be used with community data (except maybe when beta diversity is very low).

Impact on Ordination History

- by 1970 PCA was the ordination method of choice for community data
- simulation studies by Swan (1970) & Austin & Noy-Meir (1971) demonstrated the horseshoe effect and showed that the linear assumption of PCA was not compatible with the nonlinear structure of community data
- stimulated the quest for more appropriate ordination methods.