PLSC 597: Modern Measurement

Scaling

February 1, 2018

The Plan

- Scaling Overview
- Unidimensional Scaling
- Scale Reliability
- Multidimensional Scaling

Some Terms

- Always $i \in \{1, 2, ...N\}$ observations on $k \in \{1, 2, ...K\}$ indicators.
- Z or $Z_1, Z_2, ...$ will indicate the underlying / latent trait(s) / phenomena
- $D_1, D_2, ...D_K$ are dichotomous indicators
- $Y_1, Y_2, ... Y_K$ are *continuous* indicators

Scaling, In General

Characteristics:

- Comparative vs. Non-Comparative Scaling
- Subject-centered vs. Stumulus-centered
- Metric vs. Non-Metric
- Unidimensional vs. Multidimensional

(Unidimensional) Scaling: History

Thurstone (1927, 1929)

- Non-comparative, subject-centered
- "Law of Comparative Judgment": The degree to which any two stimuli can be discriminated is a direct function of the difference in their status as regards the attribute in question.
- Methods: paired comparisons, successive intervals, and equal-appearing intervals.

Likert (1932)

- Non-comparative, subject-centered
- ullet Summative + unidimensional o item construction & selection are key

Gutman (1944, 1950) ("scalogram analysis")

- · Comparative; both subject- and stimulus-centered
- The response to each item is a simple function of the sum score

See McIver and Carmines (1981) for more...

Scaling: Dissimilarities and Distances

Dissimilarities Matrix:

$$\Delta = \begin{bmatrix} 0 & \delta_{21} & \delta_{31} & \dots & \delta_{K1} \\ \delta_{12} & 0 & \delta_{32} & \dots & \delta_{K2} \\ \delta_{13} & \delta_{23} & 0 & \dots & \delta_{K3} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ \delta_{1K} & \delta_{2K} & \delta_{3K} & \dots & 0 \end{bmatrix}$$

where

$$\delta_{ij} = f(\mathbf{Y}_i - \mathbf{Y}_j)$$

Distance Examples

Euclidean ("L2") Distance:

$$\delta_{ij} = \sqrt{(Y_{i1} - Y_{j1})^2 + (Y_{i2} - Y_{j2})^2 + ...(Y_{iN} - Y_{jN})^2}$$

Manhattan ("L1") Distance:

$$\delta_{ij} = |Y_{i1} - Y_{j1}| + |Y_{i2} - Y_{j2}| + ... |Y_{iN} - Y_{jN}|$$

Minkowski Order-p ("LP") Distance:

$$\delta_{ij} = \left(\sum_{i=1}^{n} |x_i - y_i|^p\right)^{1/p}$$

Goal: Location N points in a low (p)-dimensional space such that the Euclidean distances between them approximate Δ .

That is, find a set of $N \times P$ points **D** such that $d_{ij}(\mathbf{D}) \approx \delta_{ij} \ \forall i, j$, where

$$d_{ij}(\mathbf{D}) = \sqrt{\sum_{p=1}^{P} (x_{ip} - x_{jp})^2}$$

and $p = \{1, 2, ... P\}$ denotes the dimensionality of the space.

For a $K \times K$ dissimilarity matrix Δ , the *stress* associated with a given set of coordinates D is:

$$\sigma(\mathbf{D}) = \sum_{i < j} w_{ij} [\delta_{ij} - d_{ij}(\mathbf{D})]^2$$

with weights

$$\mathbf{W} = \begin{bmatrix} 0 & w_{21} & w_{31} & \dots & w_{K1} \\ w_{12} & 0 & w_{32} & \dots & w_{K2} \\ w_{13} & w_{23} & 0 & \dots & w_{K3} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ w_{1K} & w_{2K} & w_{3K} & \dots & 0 \end{bmatrix}$$

such that

$$\sum_{i < j} w_{ij} \delta_{ij}^2 = \frac{N(N-1)}{2}.$$

Scaling Types

Key: Transformations of the dissimilarities.

Ratio:

- \cdot $\hat{d}_{ij} = \delta_{ij}$ (no transformation)
- · A special case of metric scaling

Interval:

- $\cdot \ \hat{d}_{ij} = a + b(\delta_{ij})$ (linear transformation)
- · Also *metric*; "the ratio of differences of distances should be equal to the corresponding ratio of differences in the data"

Nonlinear:

- $\cdot \hat{d}_{ij} = g(\delta_{ij})$ (e.g., splines)
- · Also metric

Ordinal:

- $\hat{d}_{ij} = f(\delta_{ij})$ such that $\delta_{ij} < \delta_{i'j'} \Rightarrow f(\delta_{ij}) < f(\delta_{i'j'})$
- · Monotone / rank-preserving transformations
- · Nonmetric

Unidimensional Scaling: The Sum Score

Simplest approach: the sum score:

$$\hat{Z}_i = \sum_{k=1}^K Y_{ik}$$

or

$$\hat{Z}_i = \frac{\sum_{k=1}^{K_i} Y_{ik}}{K_i}.$$

Requires:

- $Var(Y_j) = Var(Y_k) \forall j \neq k$
- $Cov(Y_j, Z) = Cov(Y_k, Z) \forall j \neq k$

Unidimensional Scaling

For p = 1, the stress

$$\sigma(\mathbf{D}) = \sum_{i < i} w_{ij} [\delta_{ij} - d_{ij}(\mathbf{D})]^2$$

can be shown to be a function of the orders of the rank orders of the items (see Mair and deLeeuw 2014 for details). This means:

- No need to chose among distances / similarities.
- Solution is via combinatorics.

Example: Cities

> Cities[,c(1,4)]

```
city longitude
New York
                New York -74.00594
Los Angeles Los Angeles -118.24368
Chicago
                 Chicago -87.62980
Houston
                 Houston -95.36980
Philadelphia Philadelphia -75.16522
Phoenix
                 Phoenix -112.07404
San Antonio
             San Antonio -98.49363
               San Diego -117.16108
San Diego
Dallas
                  Dallas -96.79699
                San Jose -121.88633
San Jose
```

Cities: Euclidean Distance Matrix (Δ)

```
> CityLong <- data.frame(t(Cities$longitude)) # longitudes in a row
> colnames(CityLong) <- t(Cities$city) # names
> D1long <- dist(t(CityLong)) # distance object
> D1long
             New York Los Angeles Chicago Houston Philadelphia Phoenix San Antonio San Diego Dallas
Los Angeles
                 44.2
Chicago
                 13.6
                             30.6
                 21.4
                             22.9
                                      7.7
Houston
Philadelphia
                 1.2
                             43.1
                                     12.5
                                             20.2
Phoenix
                 38.1
                              6.2
                                     24.4
                                             16.7
                                                          36.9
San Antonio
                24.5
                             19.8
                                     10.9
                                              3.1
                                                          23.3
                                                                  13.6
San Diego
                 43.2
                             1.1
                                     29.5
                                             21.8
                                                          42.0
                                                                   5.1
                                                                              18.7
Dallas
                 22.8
                             21.4
                                      9.2
                                             1.4
                                                          21.6
                                                                  15.3
                                                                               1.7
                                                                                        20.4
```

26.5

46.7

9.8

23.4

4.7

25.1

San Jose

47.9

3.6

34.3

Unidimensional Scaling

```
> library(smacof)
> UDS <- uniscale(D1long)
> UDS
Call: uniscale(delta = D1long)
Final stress value: 4.1e-16
Number of accepted permutations: 180836
Number of possible permutations: 3628800
Number of objects: 10
> UDS$conf
   New York Los Angeles
                              Chicago
                                          Houston Philadelphia
   -1.04236
                 0.75350
                             -0.48929
                                         -0.17508
                                                      -0.99530
    Phoenix San Antonio
                            San Diego
                                           Dallas San Jose
    0.50304
                -0.04827
                              0.70955
                                         -0.11715
                                                      0.90137
```

Cities: UDS Coordinates

East-West Locations for the Ten Largest U.S. Cities via UDS

PRINA Verkhia	Chicago	Houston Dallas San Antonio	Phoenix	Eds Angeles	San Jose

Another Example: SCOTUS Votes

> head(SCOTUS)

id	Rehnquist	Stevens	OConnor	Scalia	Kennedy	Souter	Thomas	Ginsburg	Breyer
1	0	1	0	0	0	0	0	0	0
2	0	0	0	0	0	0	0	0	0
3	0	0	0	0	0	0	0	0	0
4	1	1	1	1	1	1	1	1	1
5	0	1	0	0	1	1	0	1	1
7	1	1	1	0	1	1	0	1	1

SCOTUS: Sum Scores and Distances

```
> SumScores <- colSums(SCOTUS[,2:10],na.rm=TRUE) / nrow(SCOTUS)
> SumScores
Rehnquist Stevens
                    OConnor
                              Scalia
                                      Kennedy
                                                 Souter
                                                          Thomas Ginsburg Brever
  0.2838
           0.6906
                     0.4023
                              0.2648
                                       0.3672
                                                 0.6094
                                                          0.2451
                                                                   0.6130 0.5772
> D1SCOTUS <- dist(t(SCOTUS[,2:10]))
> D1SCOTUS
        Rehnquist Stevens OConnor Scalia Kennedy Souter Thomas Ginsburg
Stevens
            25.61
           15.94
OConnor
                   22.32
Scalia
           13.49
                   26.80
                         18.38
Kennedy
           13.27
                   23.62
                         16.31 15.62
            22.52 15.39
Souter
                         18 73 23 56
                                        20.57
Thomas
           13.45 26.81
                         18.36 10.15
                                       16.03 23.96
Ginsburg
           22.72 15.10 19.65 24.00
                                       20.88 11.27
                                                     24.64
            22.29 16.09 18.25 24.02
                                       20.66 13.42 24.78
                                                            12.69
Brever
```

SCOTUS: UDS

- > SCOTUS.UDS <- uniscale(D1SCOTUS)</pre>
- > SCOTUS.UDS

Call: uniscale(delta = D1SCOTUS)

Final stress value: 0.317

Number of accepted permutations: 347136 Number of possible permutations: 362880

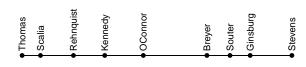
Number of objects: 9

Sum Scores and UDS

Sum Scores



UDS Results



Reliability: Cronbach's α

For a scale S that is a sum of K separate items $Y_1, Y_2, ... Y_K$,

$$\alpha = \frac{K}{K - 1} \left(1 - \frac{\sum_{k=1}^{K} \sigma_{Y_k}^2}{\sigma_{S}^2} \right)$$

where

- $\sigma_{Y_k}^2$ is the variance of item k and
- σ_S^2 is the variance of the scale S.

lpha: "The expected correlation of two tests that measure the same construct."

Cronbach's α (continued)

 α :

- "The expected correlation of two tests that measure the same construct."
- A lower bound to reliability.
- If $S = D_1 + D_2 + ... + D_K$, then $\sigma_{Y_k}^2 = P_k(1 P_k)$.
- $\alpha \in [0,1]$ (theoretically)
- Rule of thumb: $\alpha \geq$ 0.7 is "adequate"

Limitations (from Sijtsma 2009):

- Requires equal item variances, equal item covariances, and unidimensionality
- Among the "lower bounds," it's among the smallest
- A better one is the "greatest lower bound" (glb), but even it has problems...

Alpha: A Simulation

```
> # Z is the underlying thing we're trying to measure...
> N <- 400
> set.seed(7222009)
> Z < - rt(N.5)
> plot(density(Z))
> # Binary indicators... three strongly correlated with Z:
>
> Z1 <- rbinom(N,1,pbinom(Z,1,0.5))
> Z2 <- rbinom(N,1,pbinom(Z,1,0.5))
> Z3 <- rbinom(N,1,pbinom(Z,1,0.5))
>
> # ...two correlated with each other, uncorrelated with Z:
>
> Y < - rt(N.5)
> Y1 <- rbinom(N,1,pbinom(Y,1,0.5))
> Y2 <- rbinom(N,1,pbinom(Y,1,0.5))</pre>
>
> # ... and two uncorrelated with anything:
> X1 <- rbinom(N.1.0.5)
> X2 < - rbinom(N,1,0.5)
```

Alpha: A Simulation

```
> simAlpha <- alpha(simDF, check.kevs=TRUE)
Warning message:
In alpha(simDF, check.keys = TRUE) :
 Some items were negatively correlated with total scale and were automatically reversed.
This is indicated by a negative sign for the variable name.
> simAlpha
Reliability analysis
Call: alpha(x = simDF, check.keys = TRUE)
 raw_alpha std.alpha G6(smc) average_r S/N ase mean sd
     0.54
             0.55 0.66
                             0.15 1.2 0.036 0.46 0.25
lower alpha upper
                  95% confidence boundaries
0.47 0.54 0.61
Reliability if an item is dropped:
   raw_alpha std.alpha G6(smc) average_r S/N alpha se
       0.39
               0.40 0.53
                               0.10 0.68 0.047
7.1
              0.43 0.55 0.11 0.75 0.045
7.2
       0.42
Z3
     0.41
              0.42 0.56 0.11 0.73 0.046
     0.50
              0.51 0.58 0.15 1.04 0.038
Y1-
Y2- 0.49
            0.51 0.57 0.15 1.03 0.039
X1- 0.61 0.61 0.72 0.21 1.59 0.031
     0.60 0.60 0.71
X2
                               0.20 1.53
                                          0.032
 Item statistics
     n raw.r std.r r.cor r.drop mean
Z1 400 0.70 0.71 0.711 0.511 0.33 0.47
Z2 400 0.65 0.66 0.657 0.450 0.34 0.48
```

Z3 400 0.66 0.68 0.654 0.470 0.31 0.46 Y1- 400 0.52 0.51 0.467 0.270 0.62 0.49 Y2- 400 0.52 0.52 0.468 0.277 0.62 0.49 X1- 400 0.27 0.26 -0.014 -0.019 0.50 0.50 X2 400 0.30 0.29 0.020 0.013 0.51 0.50

Alpha: A Simulation

```
> # Now with only the items that "should" be there:
>
> simAlpha2 <- alpha(simDF[,1:3], check.keys=TRUE)
> simAlpha2
Reliability analysis
Call: alpha(x = simDF[, 1:3], check.keys = TRUE)
 raw_alpha std.alpha G6(smc) average_r S/N ase mean sd
    0.84
         0.84
                   lower alpha upper 95% confidence boundaries
0.82 0.84 0.87
Reliability if an item is dropped:
  raw_alpha std.alpha G6(smc) average_r S/N alpha se
Z1 0.77 0.77 0.63 0.63 3.4 0.023
Z2 0.77 0.77 0.63 0.63 3.4 0.023
Z3 0.80 0.80 0.67 0.67 4.1 0.020
```

Item statistics

n raw.r std.r r.cor r.drop mean sd Z1 400 0.88 0.88 0.79 0.72 0.33 0.47 Z2 400 0.88 0.88 0.79 0.72 0.34 0.48 Z3 400 0.86 0.86 0.75 0.69 0.31 0.46

Reliability: SCOTUS Data

```
> SCOTUSAlpha <- alpha(SCOTUS[.2:10],check.kevs=TRUE)
> SCOTUSAlpha
Reliability analysis
Call: alpha(x = SCOTUS[, 2:10], check.keys = TRUE)
 raw_alpha std.alpha G6(smc) average_r S/N
                                         ase mean
                                                   sd
      0.9
               0.9
                     0.93
                              0.49 8.6 0.0044 0.45 0.35
lower alpha upper
                   95% confidence boundaries
0.89 0.9 0.9
Reliability if an item is dropped:
        raw_alpha std.alpha G6(smc) average_r S/N alpha se
Rehnquist
             0.88
                      0.88
                             0.91
                                      0.48 7.4 0.0049
Stevens
             0.89
                      0.89
                           0.92
                                      0.51 8.4 0.0045
                          0.92
OConnor
             0.88
                      0.88
                                      0.48 7.3 0.0051
                      0.89 0.91
Scalia
           0.89
                                      0.50 7.9 0.0046
Kennedy
           0.88
                      0.88
                           0.92
                                      0.48 7.3 0.0051
                           0.91
Souter
           0.88
                      0.88
                                      0.47 7.2 0.0051
Thomas
           0.89
                      0.89
                           0.91
                                      0.50 8.0 0.0046
Ginsburg
           0.88
                      0.88
                           0.91
                                      0.48 7.4 0.0051
Breyer
            0.88
                      0.88
                           0.91
                                      0.49 7.5 0.0050
```

Multidimensional Scaling

Types...

- **Metric** Inputs are interval/ratio-level measures; transformations from δ_{ij} to d_{ij} are cardinal-valued.
- **Non-Metric** Inputs are binary or ordinal-level; transformations are rank-preserving *only*.

General steps:

- Generate a dissimilarity matrix Δ (chosing distance metric)
- Choose the dimensionality *p*
- Choose the type of scaling
- Fit model + interpret the results
- Assess model fit + conduct diagnostics

MDS: Model Fit / Diagnostics

Kruskal's rule of thumb for stress:

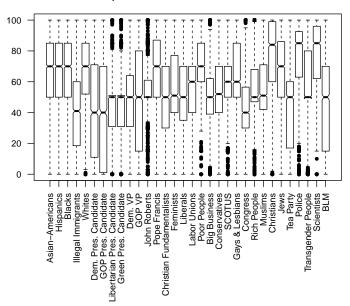
- $\sigma(\mathbf{D}) = 0.20 \rightarrow \text{"poor"}$
- $\sigma(\mathbf{D}) = 0.10 \rightarrow$ "fair"
- $\sigma(\mathbf{D}) = 0.05 \rightarrow \text{"good"}$
- $\sigma(\mathbf{D}) = 0.025 \rightarrow$ "excellent"
- $\sigma(\mathbf{D}) = 0 \rightarrow$ "perfect"

Key diagnostic: **Shepard plot**: a scatterplot of δ_{ij} vs. \hat{d}_{ij} ...

- Illustrates model "fit"
- Also illustrates the transformation of the δ_{ij} s

Also permutation tests for model fit...

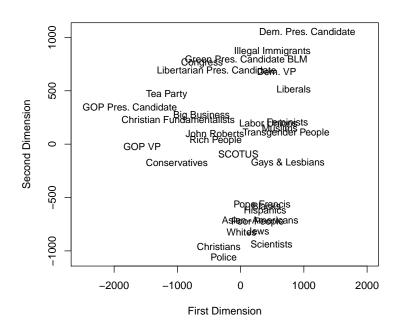
Example: 2016 ANES Thermometer Scores



MDS using cmdscale

```
> MDS2.alt <- cmdscale(ThermDist,k=2)</pre>
> head(MDS2.alt)
                            \lceil .1 \rceil \quad \lceil .2 \rceil
Asian-Americans
                         309.90 -709.9
                         387.51 -627.6
Hispanics
Blacks
                         407.03 -580.7
                         501.74 866.3
Illegal Immigrants
Whites
                          14.72 -821.5
Dem. Pres. Candidate 1054.53 1053.2
```

MDS Plot, using cmdscale



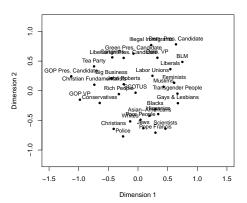
MDS using mds (ratio scaling)

```
> MDS2 <- mds(ThermDist, ndim=2)
> MDS2
```

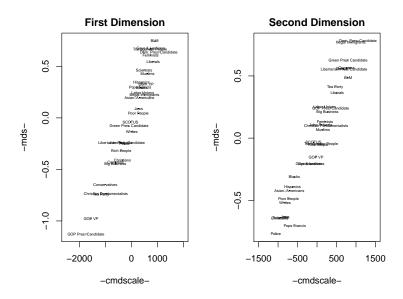
Call:

mds(delta = ThermDist, ndim = 2)

Model: Symmetric SMACOF Number of objects: 32 Stress-1 value: 0.221 Number of iterations: 87



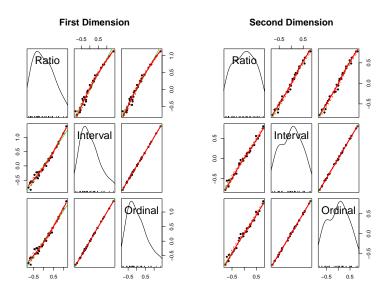
cmdscale and mds Comparison



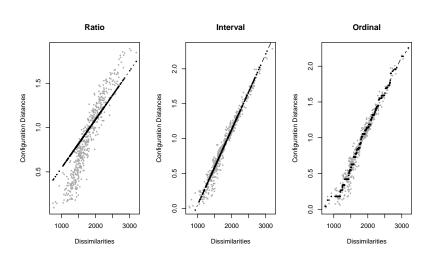
Alternatives: Interval and Ratio Scaling

```
> MDS2.int <- mds(ThermDist, ndim=2, type="interval")
> MDS2 int
Call:
mds(delta = ThermDist, ndim = 2, type = "interval")
Model: Symmetric SMACOF
Number of objects: 32
Stress-1 value: 0.087
Number of iterations: 17
> # Ordinal:
> MDS2.ord <- mds(ThermDist, ndim=2, type="ordinal")
> MDS2 ord
Call:
mds(delta = ThermDist, ndim = 2, type = "ordinal")
Model: Symmetric SMACOF
Number of objects: 32
Stress-1 value: 0.071
Number of iterations: 22
```

Ratio-Interval-Ordinal Comparison



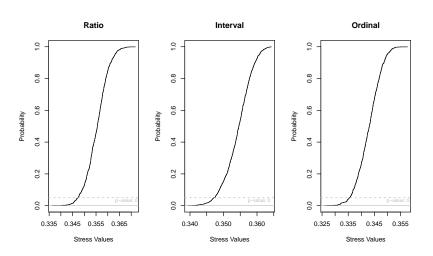
Diagnostics: Shepard Diagrams



Diagnostics: Permutation Tests

```
> ThermR.permtest<-permtest(MDS2.nrep=1000.verbose=FALSE)
> ThermR.permtest
Call: permtest.smacof(object = MDS2, nrep = 1000, verbose = FALSE)
SMACOF Permutation Test
Number of objects: 32
Number of replications (permutations): 1000
Observed stress value: 0.221
p-value: 0
> ThermI.permtest<-permtest(MDS2.int,nrep=1000,verbose=FALSE)
> ThermI.permtest
Call: permtest.smacof(object = MDS2.int, nrep = 1000, verbose = FALSE)
SMACOF Permutation Test
Number of objects: 32
Number of replications (permutations): 1000
Observed stress value: 0.087
p-value: 0
> ThermO.permtest<-permtest(MDS2.ord,nrep=1000,verbose=FALSE)
> ThermO.permtest
Call: permtest.smacof(object = MDS2.ord, nrep = 1000, verbose = FALSE)
SMACOF Permutation Test
Number of objects: 32
Number of replications (permutations): 1000
Observed stress value: 0.071
p-value: 0
```

Permutation Test Plots



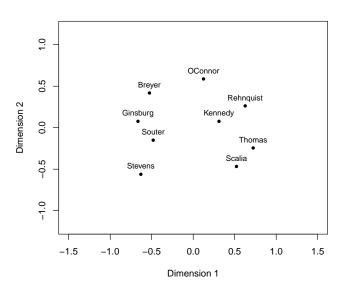
SCOTUS Redux (p = 2)

```
> SCR <- mds(D1SCOTUS, ndim=2, type="ratio")
> SCR

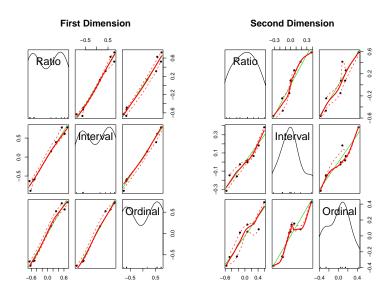
Call:
mds(delta = D1SCOTUS, ndim = 2, type = "ratio")

Model: Symmetric SMACOF
Number of objects: 9
Stress-1 value: 0.184
Number of iterations: 73
```

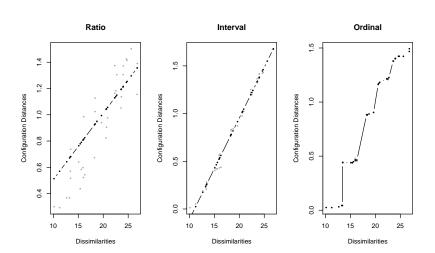
SCOTUS MDS Plot



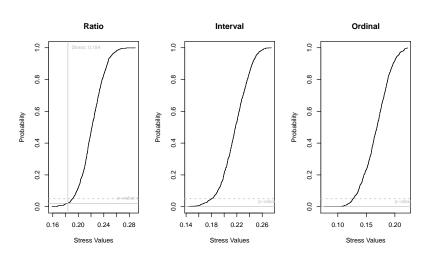
Ratio-Interval-Ordinal Comparison



SCOTUS Diagnostics: Shepard Diagrams



SCOTUS Permutation Test Plots



Useful References

- Kruskal, J.B., and M. Wish 1978. Multidimensional Scaling. Sage.
- McIver, John, and Edward C. and Carmines. 1981. *Unidimensional Scaling*. Sage Publications.
- Davison, Mark L. Multidimensional Scaling. 1983. New York: Wiley.
- Cox, Trevor F. and Michael A. A. Cox. 2000. Multidimensional Scaling, 2nd Ed. New York: Chapman & Hall.
- Borg, Ingwer, and Groenen, Patrick. 2005. Modern
 Multidimensional Scaling: Theory and Applications, 2nd Ed. Berlin:
 Springer-Verlag.
- Borg, Ingwer, Patrick Groenen, and Patrick Mair. 2013. Applied Multidimensional Scaling. Berlin: Springer-Verlag.

Useful R Packages and Routines

Distances, Proximities, etc.

- dist function (base R)
- distances package
- proxy package

Scaling

- stats::cmdscale (classical MDS, in base R)
- smacof (state-of-the-art MDS package)
- vegan (ecology package; has some good MDS routines)
- Others...

Useful Links

- smacof documentation.
- Seven ways to do MDS in R.
- Jan De Leeuw's website.