## General Crosswalk Construction Framework

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### 1 Notation formatting

- **Bold** = variable or set
- *Italic* = a single instance (= item in set)
- Non-italic = a set
- UPPER CASE = input parameter
- lower case = derived from input parameters

# 2 Goal specification

Generate a crosswalk  $\boldsymbol{X_{\mathrm{ST}}}$  ...

- To target zones **T** (geographic level  $G_{\mathbf{T}}$  in year  $Y_{\mathbf{T}}$ )
- Including exactly one record per atom st (an intersection between source zone s and target zone t)
- With interpolation weights  $w_{ST}$ 
  - A single weight  $w_{cst}$  for each count variable c in C, for each atom st
    - \*  $w_{\text{cst}} = \text{proportion of } c \text{ in } s \text{ (denominator) that is also in } st \text{ (numerator)} = \frac{c_{st}}{c_s}$
    - \* All  $\mathbf{C}$  are count variables (e.g., population, housing units, etc.) that have been reported for a set of sub-zones  $\mathbf{S}'$  (blocks)
- Build from an existing crosswalk  $X_{\mathbf{S'T'}}$  ...
  - From source sub-zones S', which nest within S
  - To source sub-zones T', which nest within T
  - In our setting, we can assume:
    - \*  $G_{S'} = G_{T'} = blocks$
    - \*  $Y_{S'} = Y_S$  and  $Y_{T'} = Y_T$
  - Includes weights  $w_{S'T'}$  indicating proportion of each source sub-zone's features (population & housing) in each sub-zone atom s't'.
- Include every s in S and every t in T.
  - Atom records may have *null* s where t where a zone in one set lies beyond the spatial extent of the other set, or the intersection is outside the extent of  $X_{S'T'}$ .
    - \* In our case,  $X_{S'T'}$  is a block-to-block crosswalk based on NHGIS shapefiles, which are clipped at the coast. The 1990-2010 crosswalk omits "off-coast" 1990 blocks that are not in the shapefile. For crosswalks with 1990 source zones, we may have  $null\ t$  for source zones that lie entirely off-coast.

# 3 Summary of key input parameters

- $G_S$  = source geographic level
- $Y_S$  = source year
- $G_{\mathbf{T}} = \text{target geographic level}$
- $Y_T = \text{target year}$
- $\bullet$  C = set of count variables for which to derive separate weights

# 4 General steps

- 1. Obtain & load sub-zone crosswalk (blocks-to-blocks)  $\boldsymbol{X_{S'T'}}.$
- 2. Obtain & load data for source sub-zone counts (source-year block data)  $C_{\mathbf{S}'}$ .
  - (a) Include any identifiers needed to associate S' with S.
- 3. Join base crosswalk  $X_{S'T'}$  to source sub-zone data  $C_{S'}$  on S' identifiers.
  - (a) Use a "left join" to ensure that all sub-zone atoms are included, even those without a matching record in the sub-zone data file (especially important for 1990 blocks).
- 4. For each sub-zone atom s't', identify encompassing zones s and t:
  - (a) If possible, derive S and T identifiers from S' and T' identifiers (e.g., tract ID is in block ID).
  - (b) Else if possible, derive **S** identifiers from source sub-zone data from step 2.
  - (c) Else, obtain identifiers through other means...
    - i. 1990 block-group parts require some special handling because neither 4a nor 4b pertain to all BGPs.
    - ii. If we generate crosswalks for target zones that cannot be identified from block IDs (e.g., places, county subdivisions, etc.), we'll need to add a step to join block crosswalk to target-year block data that includes identifiers for the target zones.
  - (d) Where s' is null (= ""), omit these dummy sub-zone atoms from subsequent computations.
    - i. This may drop some valid t from the computations, but step 9 will re-add them if needed.
- 5. Compute counts for all weighting variables in each sub-zone atom:  $c_{S'T'} = w_{S'T'} * C_{S'}$ .
- 6. Compute counts for all weighting variables in each atom of interest:  $c_{ST} = \sum c_{S'T'}$  group by S, T.
  - (a) Steps 5 & 6 can be combined into single formula by substituting  $w_{\mathbf{S'T'}} * C_{\mathbf{S'}}$  for  $c_{\mathbf{S'T'}}$  in step 6.
- 7. Compute counts for all weighting variables in each source zone:  $c_{\mathbf{S}} = \sum c_{\mathbf{ST}}$  group by **S**.
- 8. Compute all weights for all atoms of interest:  $w_{\text{CST}} = \frac{c_{st}}{c_{s}}$ .
  - (a) If  $c_s = 0$ , set  $w_{cst} = 0$ .
- 9. If  $w_{cst}$  is missing data for any s in S or t in T, add dummy atoms with null t for non-null s or null s for non-null t, and set wCst = 0.
- 10. Export clean, complete file for distribution.
  - (a) Exact specifications TBD.