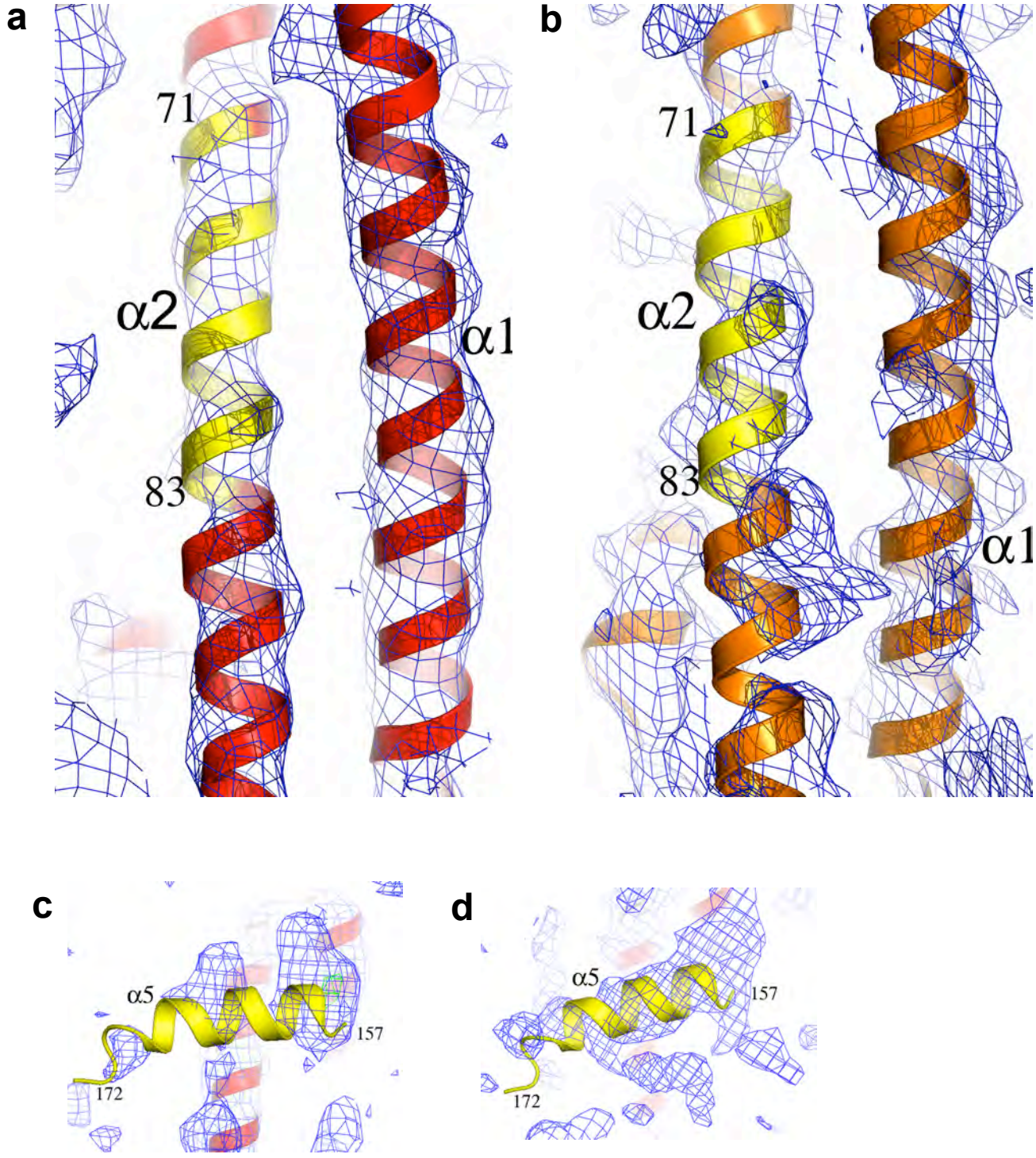

Supplementary Figure 2 Representative and unbiased electron density (2Fo-Fc, 1σ x RMSD) for the CHMP3 crystal structures. Regions in yellow were omitted from the model prior to molecular replacement.
(a) $CHMP3_{8-222}$ α1 and α2 region (residues 71-83 omitted) (4.0 Å resolution).
(b) $CHMP3_{1-150}$ α1 and α2 region (residues 71-83 omitted) (3.7Å resolution).
(c,d) Unbiased density for the autoinhibitory helix in both molecules of the $CHMP3_{8-222}$ asymmetric unit.