

**S2 Fig. Electron density maps of ligands bound in the lipid-binding site.** (A) Electron density ( $3\sigma$  level; green wires) of a fragment of 06:0 PG (sticks) in the initial difference Fourier map calculated after molecular replacement. (B)  $2F_o-F_c$  density ( $1\sigma$  level; blue wires) for the 06:0 PG fragment (sticks) after the last round of refinement. (C) Electron density ( $3\sigma$  level; green wires) of a fragment of 06:0 PA (sticks) in the initial difference Fourier map calculated after molecular replacement. (D)  $2F_o-F_c$  density ( $1\sigma$  level; blue wires) for the 06:0 PA fragment (sticks) after the last round of refinement. (E) Initial difference Fourier positive electron density ( $3\sigma$  level; green wires) of a glycerol molecule (sticks) bound in the lipid-binding site in the 10E8-T117v2 structure. (F)  $2F_o-F_c$  density ( $1\sigma$  level; blue wires) calculated at final stage of refinement for the glycerol molecule (sticks) bound in the lipid-binding site in the 10E8-T117v2 structure.

