

Intermolecular structure and hydrogen-bonding in liquid 1,2-propylene carbonate and 1,2-glycerol carbonate determined from neutron scattering

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Supplemental Information

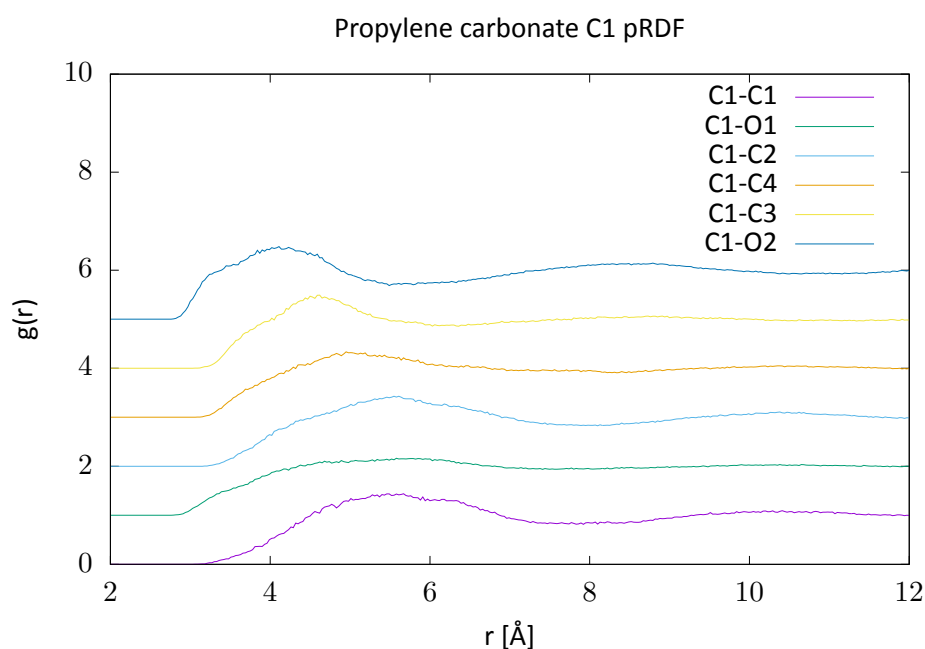


Figure S1 pRDFs to propylene carbonate C1

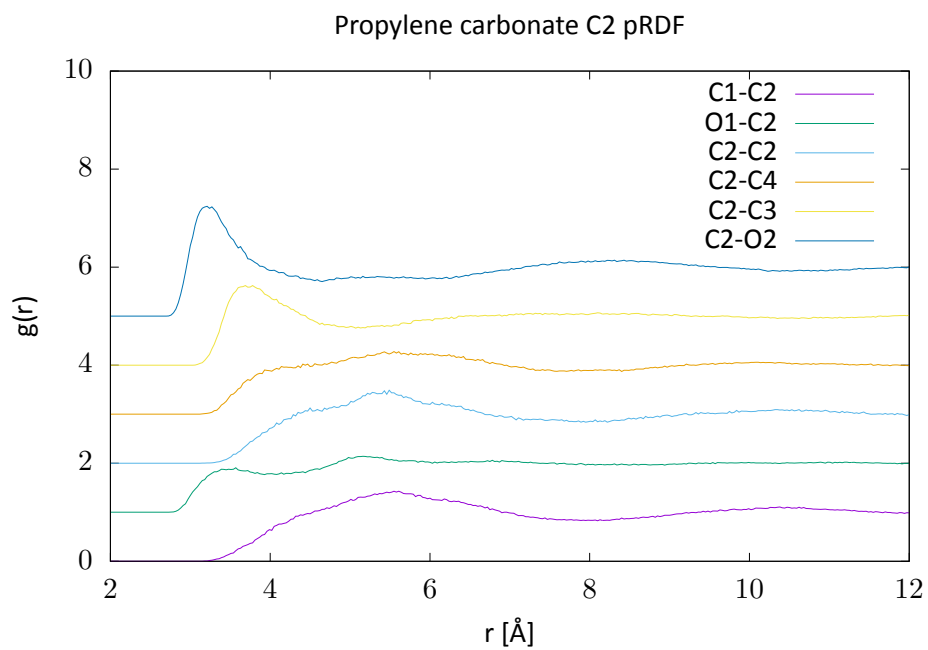


Figure S2 pRDFs to propylene carbonate C2

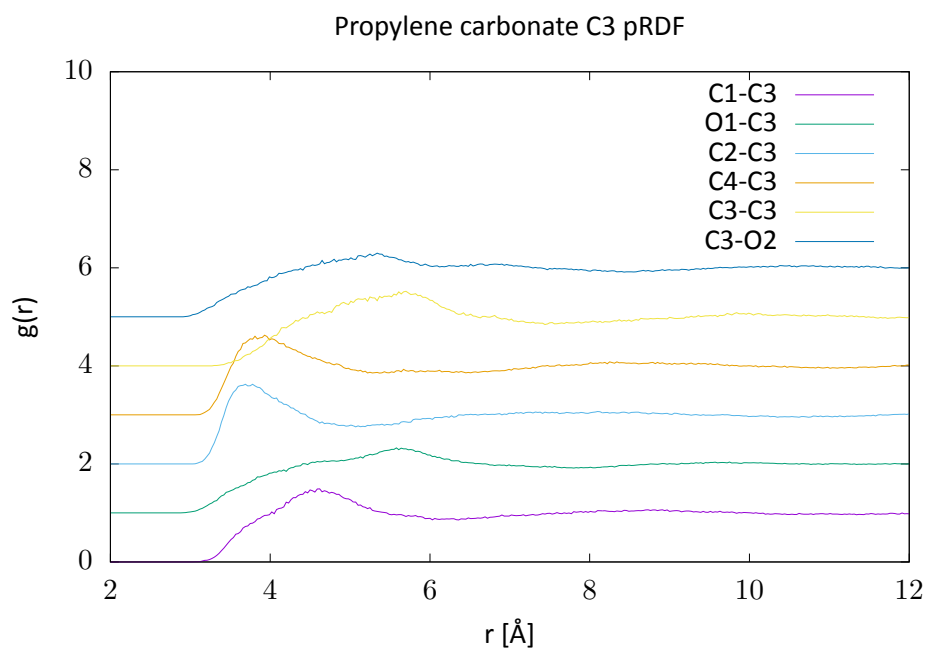


Figure S3 pRDFs to propylene carbonate C3

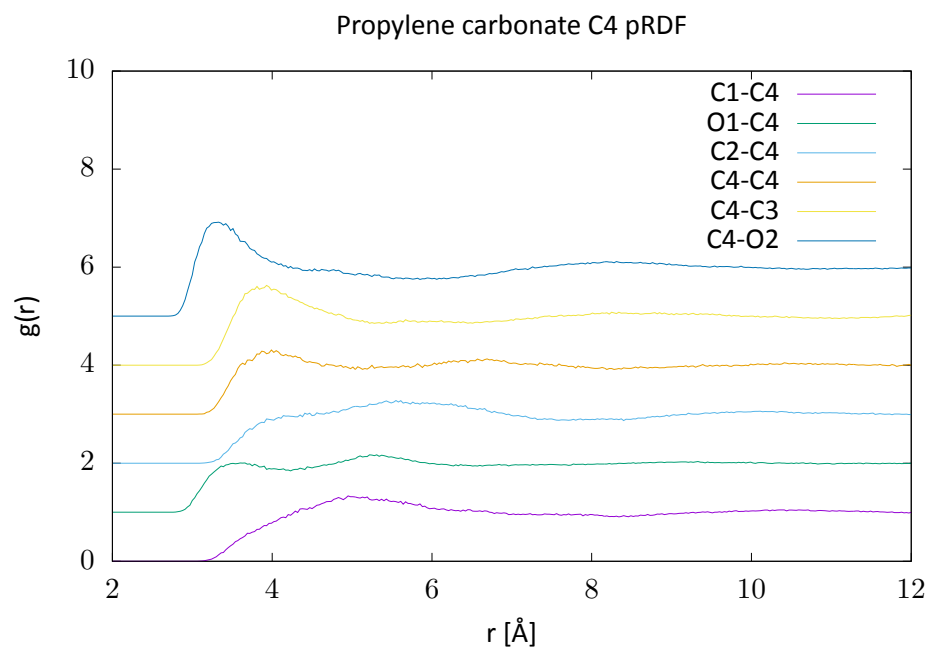


Figure S4 pRDFs to propylene carbonate C4

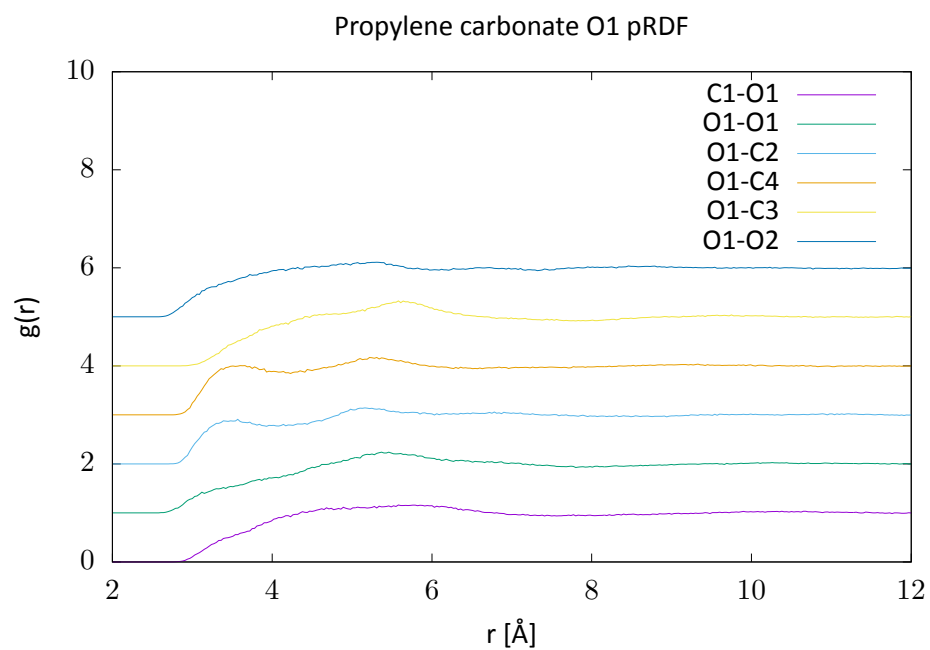


Figure S5 pRDFs to propylene carbonate O1

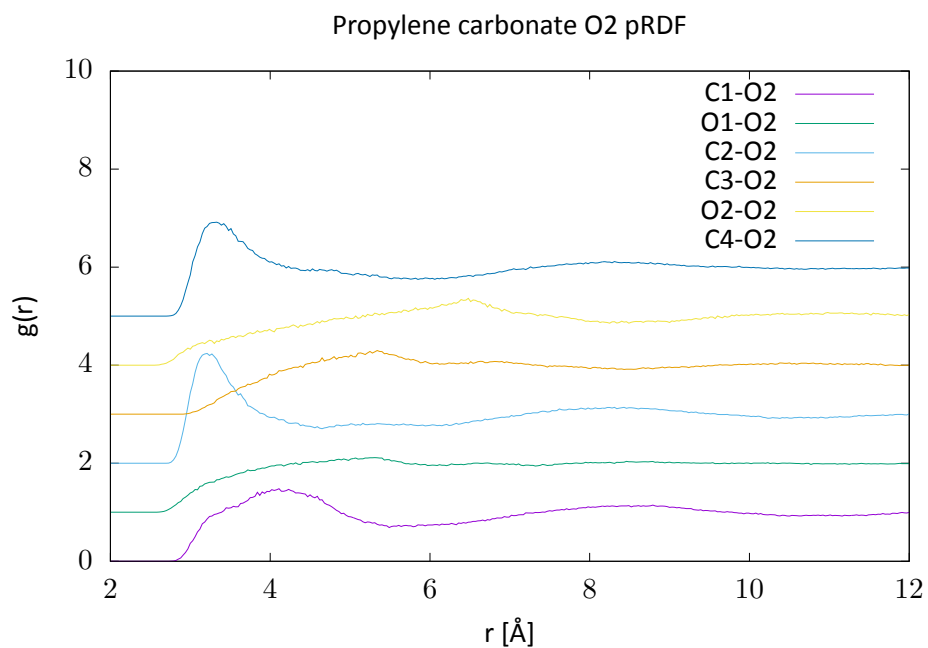


Figure S6 pRDFs to propylene carbonate O2

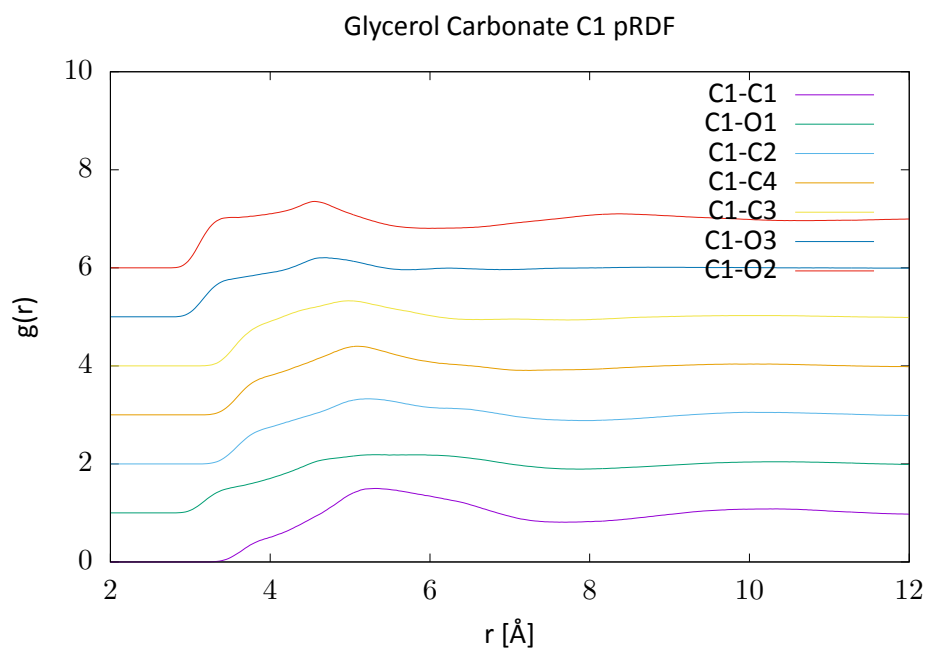


Figure S7 pRDFs to glycerol carbonate C1

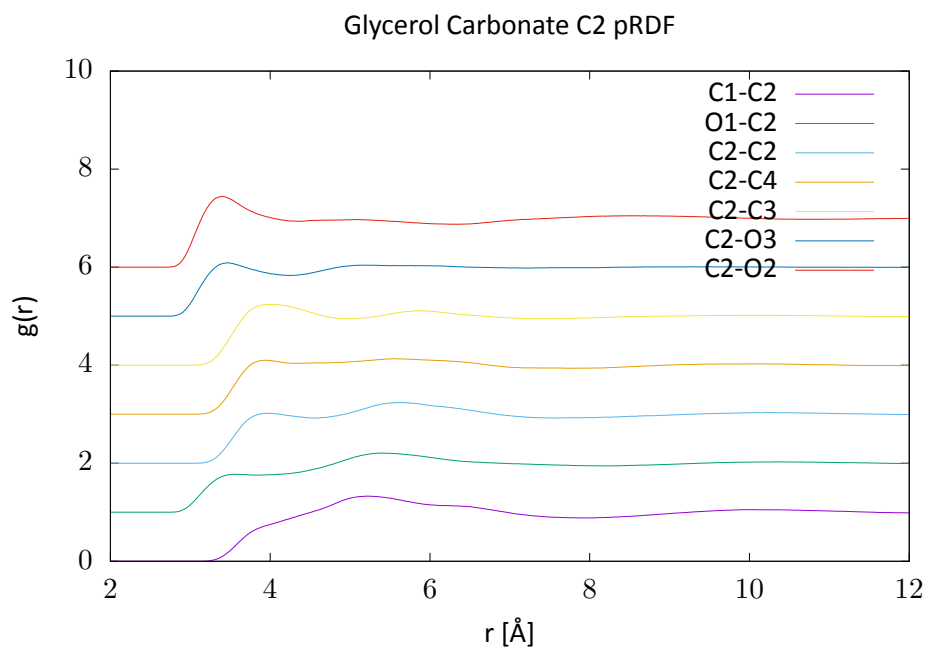


Figure S8 pRDFs to glycerol carbonate

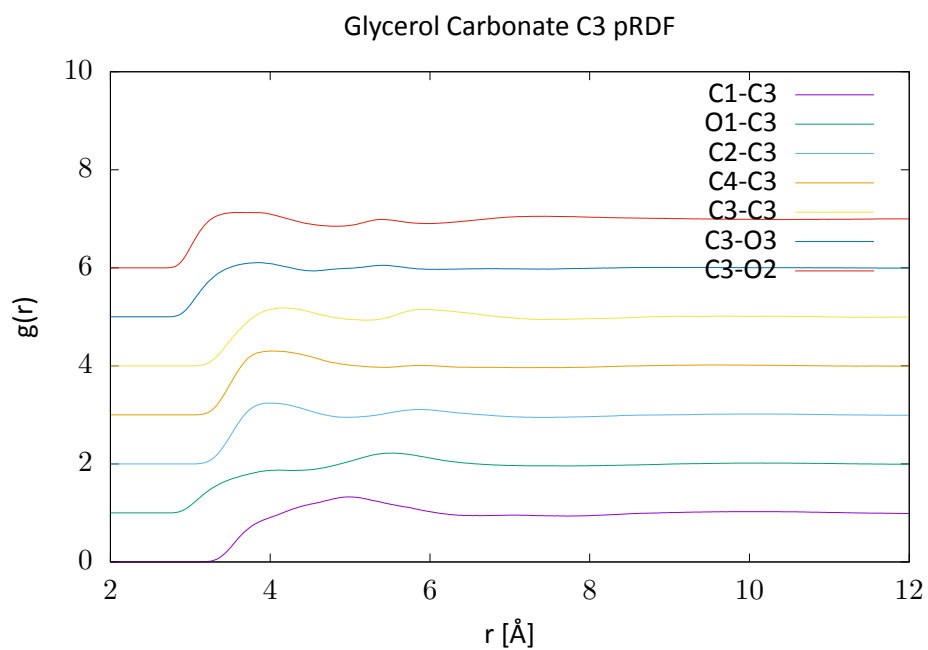


Figure S9 pRDFs to glycerol carbonate

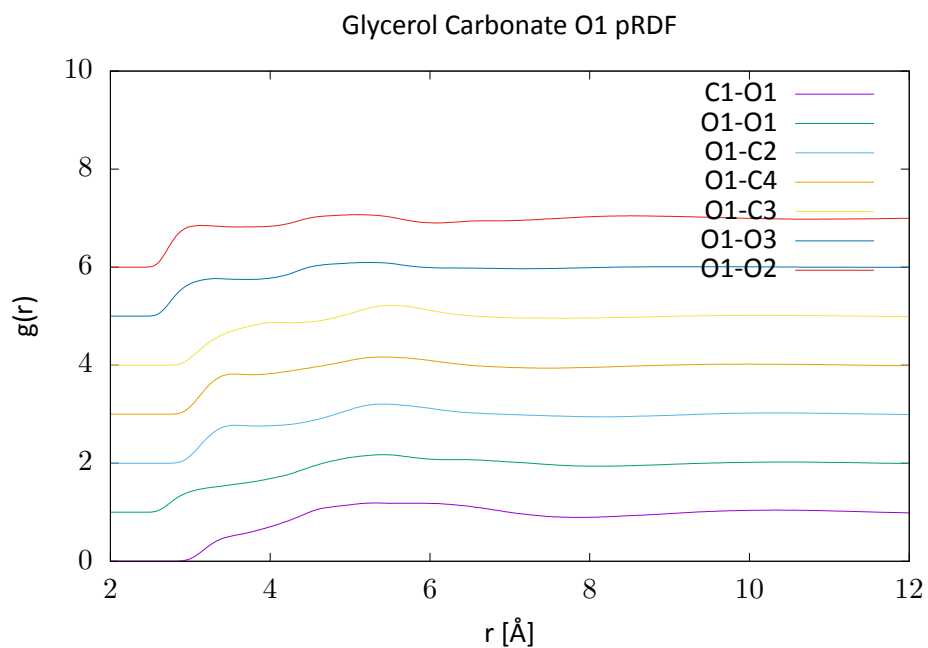


Figure S10 pRDFs to glycerol carbonate

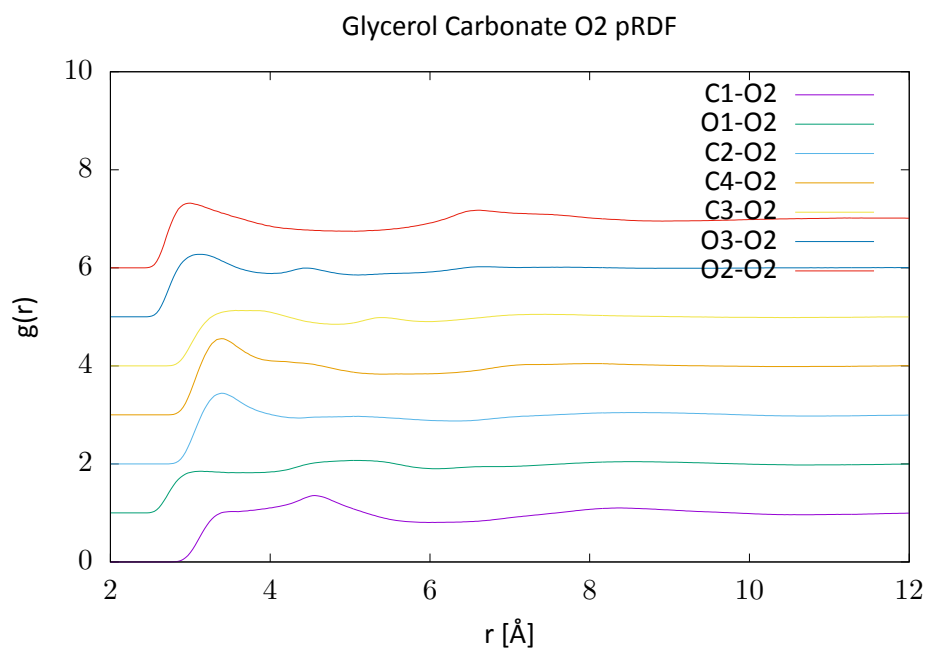


Figure S11 pRDFs to glycerol carbonate

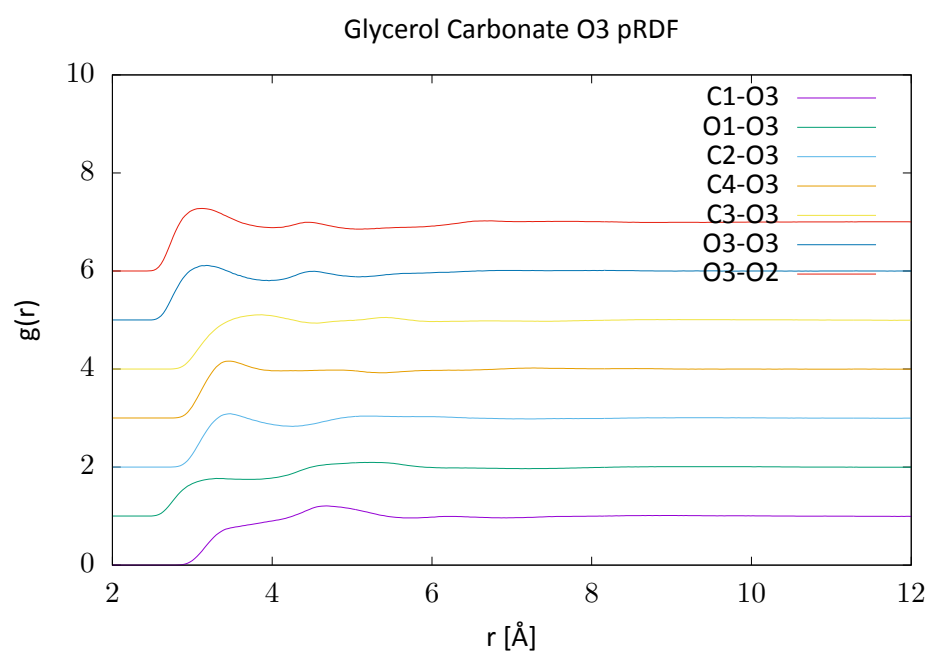


Figure S12 pRDFs to glycerol carbonate

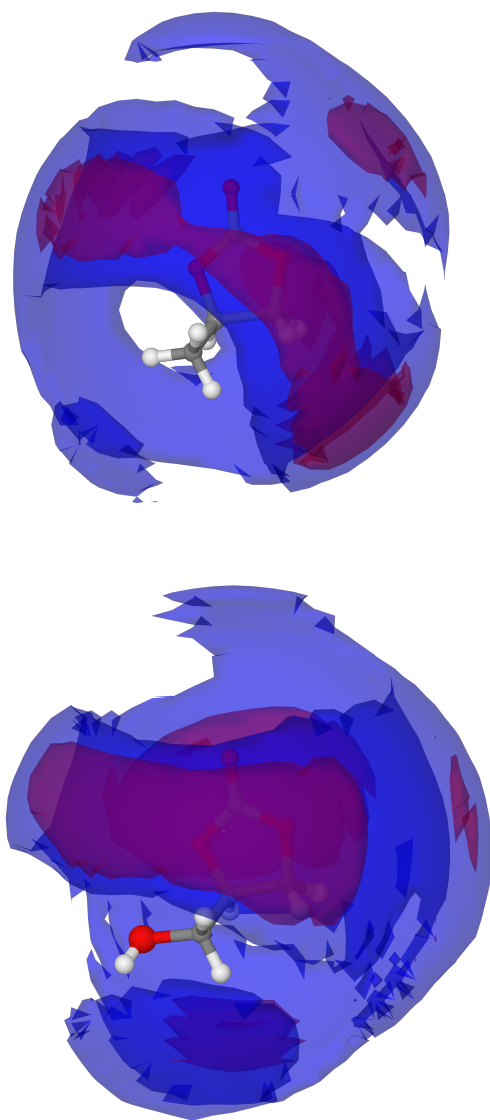


Figure S13 Centre of mass SDF volumes for propylene carbonate (top) and glycerol carbonate (bottom) calculated over the first coordination shell (3–6.5 Å) to capture the top 25 % probability density (blue) and top 5 % density (red).