

Early stages of the formation of nanocrystalline C-S-H colloids

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Calcium silicate hydrate is the primary binding component of concrete, making it one of the most important construction material of the modern world. Nevertheless, due to its complex nano-crystalline and nanocomposite nature, the large varieties of defects that can occur within its structure and the fast dissolution/reprecipitation dynamics, its formation mechanism is still a subject of debate. Previous studies revealed that calcium silicate hydrate forms through a two-stage process: i.e. a fast nucleation of nano-crystalline colloidal nuclei, followed by a slower formation of aggregates that retain its initial nano-crystalline morphology.

Here we present a new set of experiments based on state-of-the-art, in situ, High Energy X-ray Scattering (HEXS) and pair distribution function analysis, which allowed us to resolve the basic structural motifs of the initial building blocks that are formed prior to the crystallization of β -CSH. The identification of these initial building blocks will clarify the structural transitions through which CSH is formed and consequently help us understanding the basic kinetics of CSH-formation processes. A fundamental understanding of these processes will contribute towards high-performance custom-made cements.