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Modelling of degradable vinyl polymer production

Aim

Finding reactor operator conditions to tune the molecular structure of degradable vinyl polymers.

Justification

Traditional vinyl polymers are not biodegradable but display excellent material properties. In contrast, biodegradable polymers such as polyesters can be easily (bio)degraded via hydrolysis but lack high mechanical strength. The best of both worlds can be achieved by incorporating ester moieties along the vinyl backbone, providing susceptibility to hydrolysis and achieving (bio)degradation, which is an ambitious task as monomers for polycondensation and radical polymerization processes cannot react in one single process. One promising route is the radical polymerization (FRP) of traditional vinyl monomers and novel cyclic ketene acetals (CKA) or DOT monomers, the ring opening of which provides the critical (bio)degradable ester moiety during FRP (see green monomer in the Figure below).



Many chemists are investigating batch processes in which the reactivity of the CKA or DOT type is tuned based on the reactivity of the vinyl monomer (styrene, (meth)acrylates, vinyl ether, *etc.*). This chemical approach reduces composition drifting to some degree, incorporating the CKA or DOT at regular distances along the vinyl backbone, providing efficient biodegradation. However, composition drifting can be nearly completely avoided by using semi-batch feeding of reactors, in which the more reactive CKA and/or DOT monomers are slowly fed over time to a reactor already containing all the less reactive vinyl monomer, thereby greatly uniformizing the vinyl segment length and achieving optimal (bio)degradability. Such reactor operation provides opportunities for advances in polymer reaction engineering to complement advances in polymer chemistry recently achieved by research labs in Marseille and Singapore, paving the way to degradable polymer production without compromising material properties of vinyl polymers.

Program

- 1. Surveying the literature and constructing a database of reactivity ratios of vinyl-CKA pairs.
- 2. Develop a polymerization model, using in-house available commercial software, for batch and semi-batch vinyl copolymerization.
- 3. Construction of the correlations between operating conditions (batch time, feeding regime, *etc.*) and the molecular structure of polymers (vinyl segment length, composition drifting, *etc.*).
 - Powerling the best operating conditions for optimal molecular structure of degradable polymers.

