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Supporting Information for

Comparative study of the hydrogen bonding interactions between ester-functionalized/non-functionalized imidazolium-based ionic liquids and DMSO

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Fig. S1. Optimized geometries of [AOEMIM]⁺, Hydrogen bonds were denoted by dashed lines, the corresponding distances of H···O was labelled.

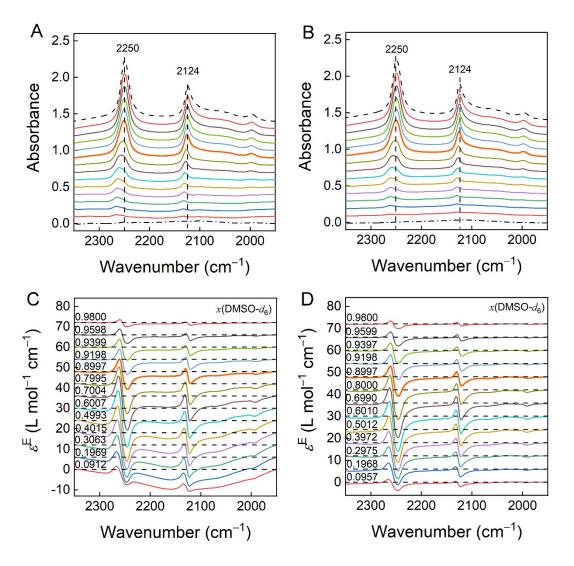


Fig. S2. ATR-FTIR and excess spectra of AOEMIMBF₄–DMSO- d_6 (A, C) and BMIMBF₄–DMSO- d_6 (B, D) systems in the range of the C–D stretching vibrations.

The dashed-dotted lines and dashed line in (A, B) depict the spectra of AOEMIMBF₄/ BMIMBF₄ and DMSO- d_6 , respectively.