Vitrification of the smectic C_A* phase and kinetics of cold crystallization investigated for the fluorinated compound with a chiral centre based on (*S*)-(+)-3-octanol

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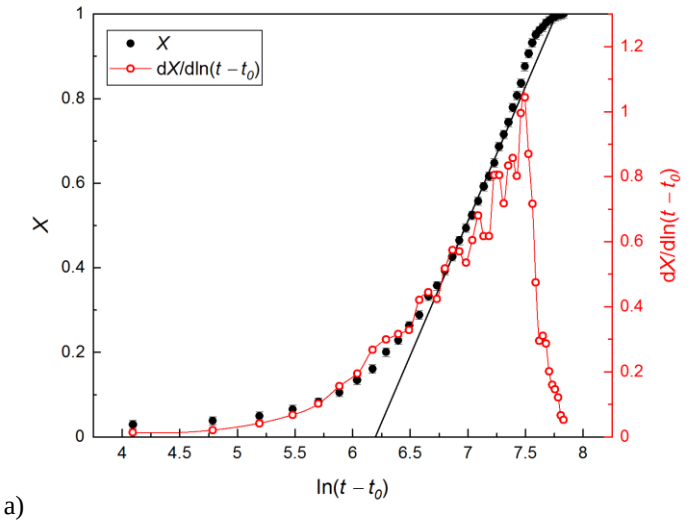
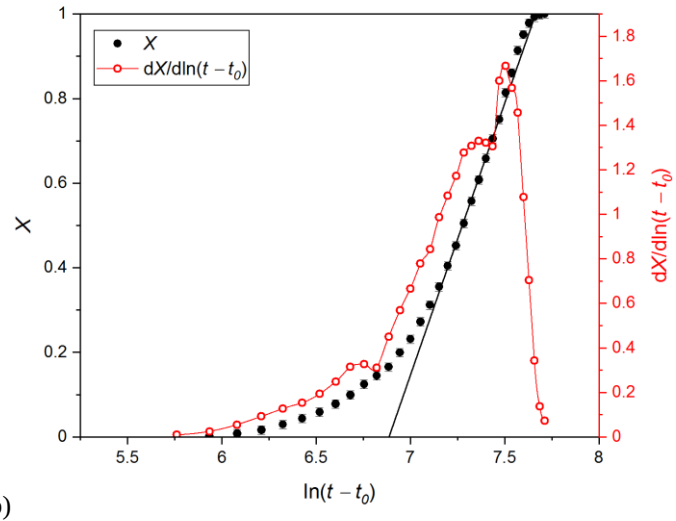
Supporting Information

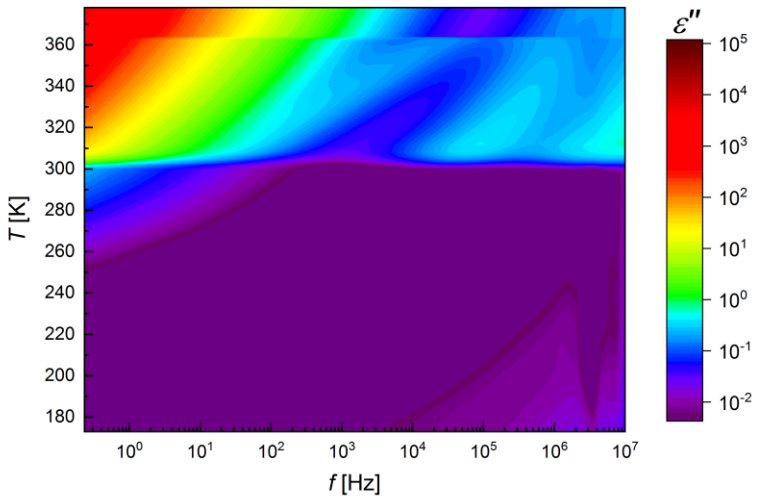


a)

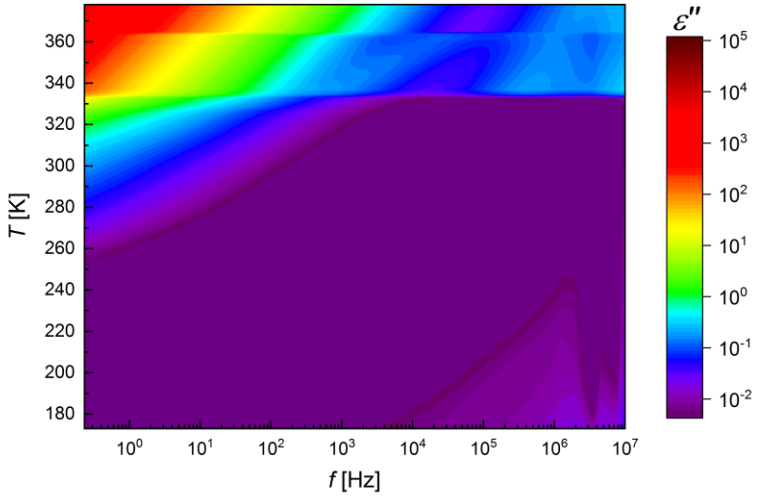


b)

Figure S1. Avramov plots for the isothermal melt crystallization of II.5.(HH) (S) in 278 K (a) and 283 K (b).



a)



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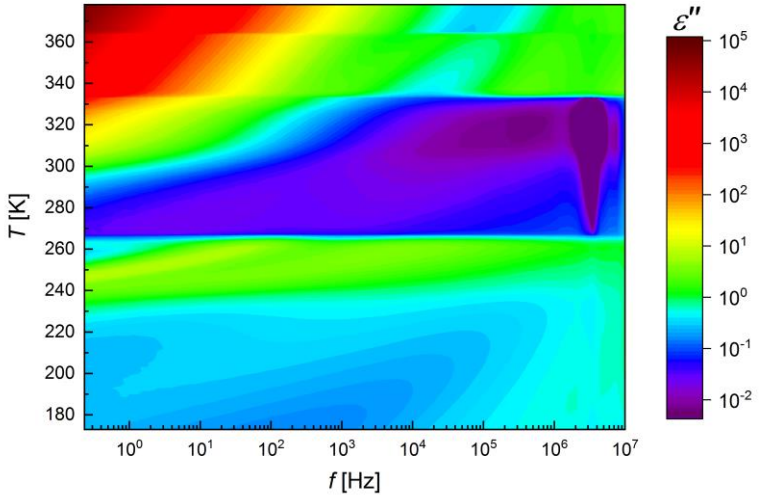
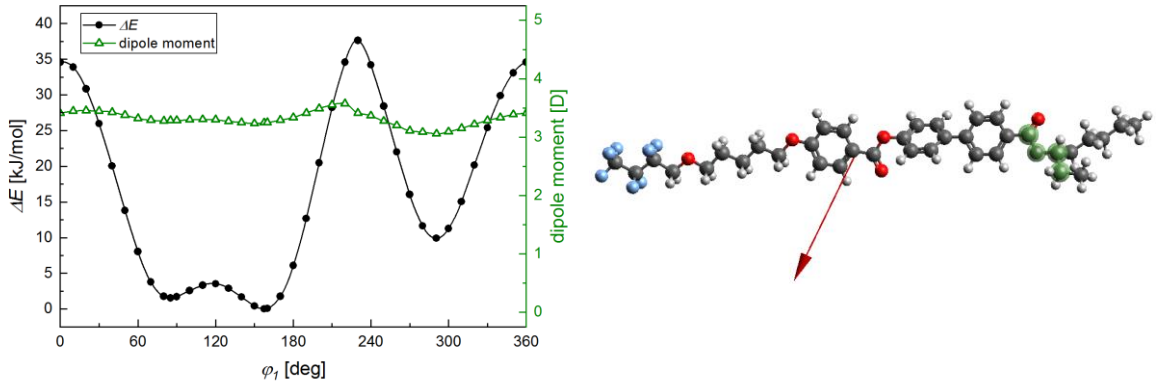
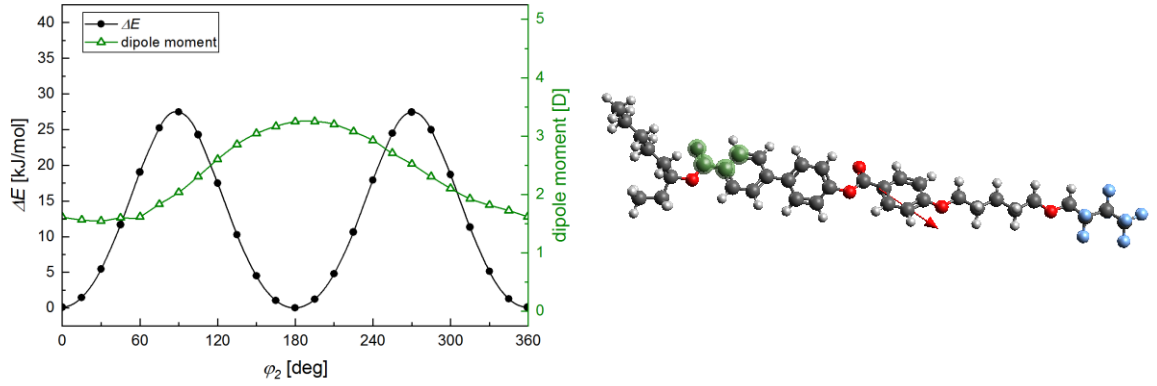


c)

Figure S2. Dielectric absorption of II.5.(HH) (S) vs. temperature and frequency on slow cooling (a), on heating after slow cooling (b) and on heating after fast cooling (c).



a)



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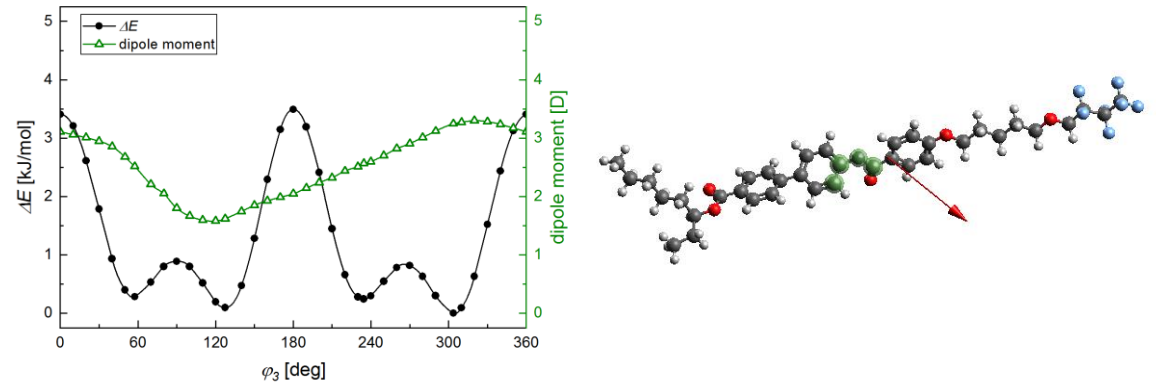
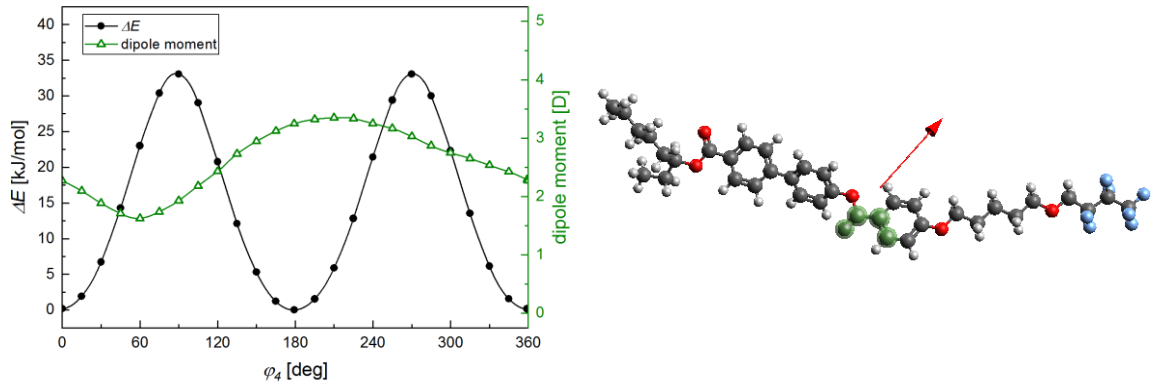
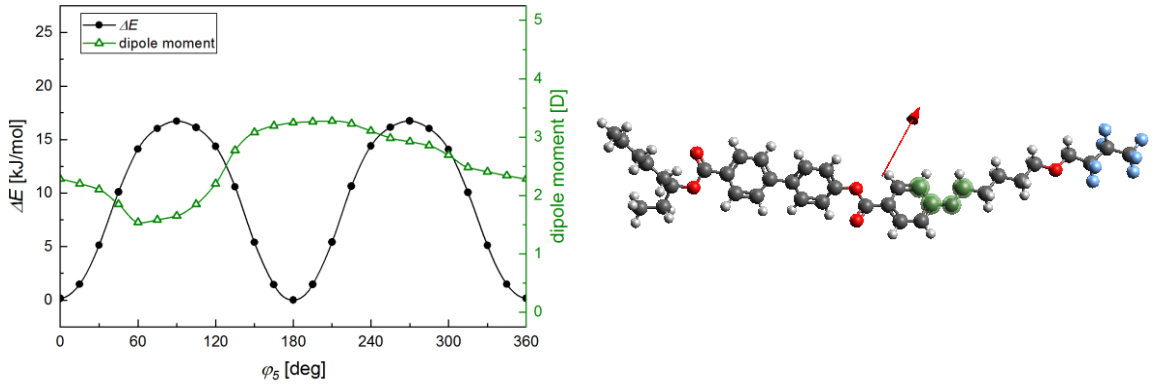


c)

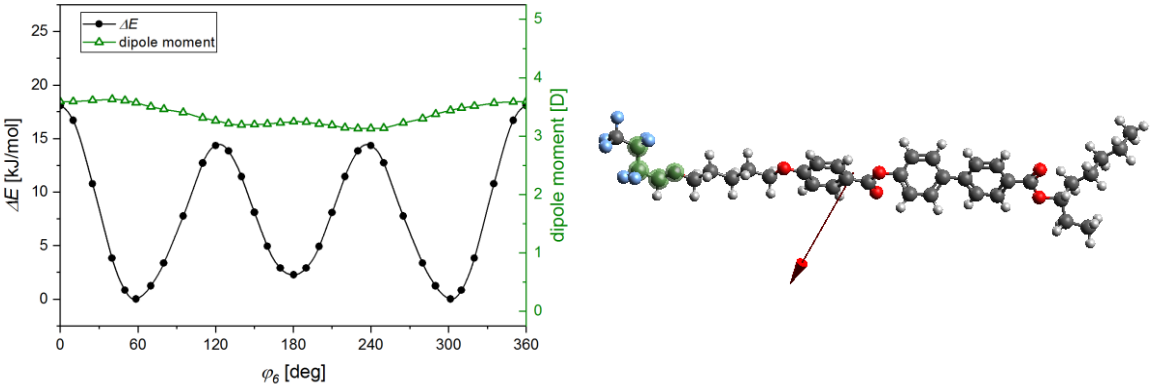


d)

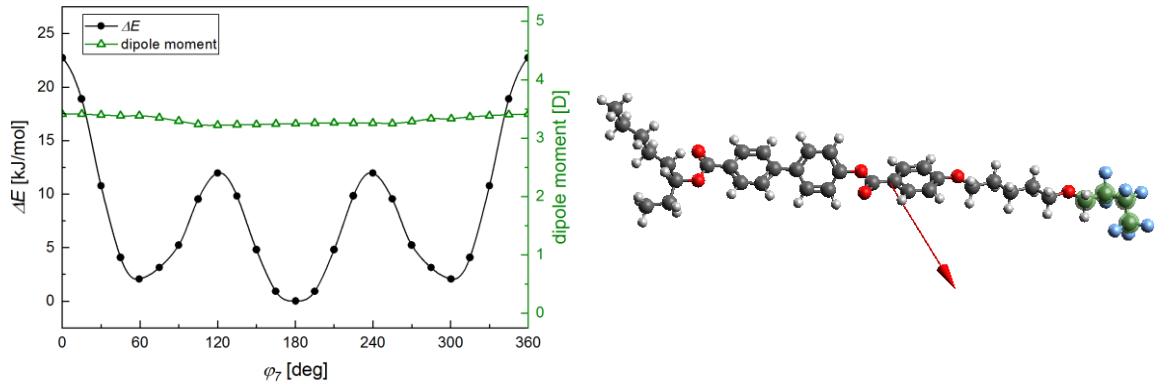
Figure S3 (part I). Conformational energy (circles, left axis) and total dipole moment (squares, right axis) vs. torsional angle for selected intra-molecular rotations calculated for the isolated II.5.(HH) (S) molecule with the DFT-B3LYP/def2TZVP method.



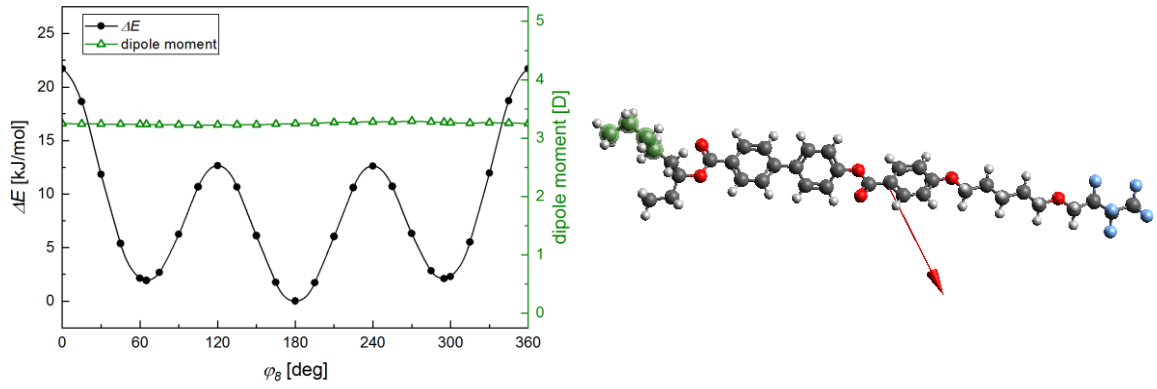
e)



f)

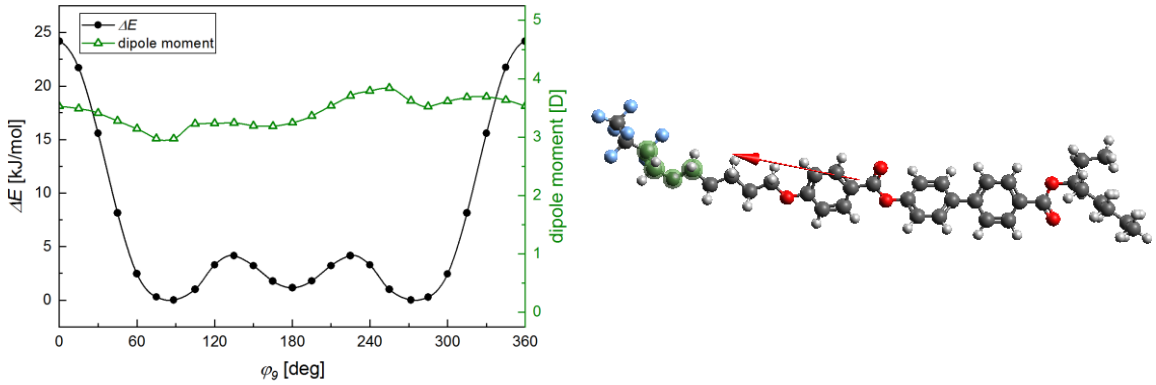


g)

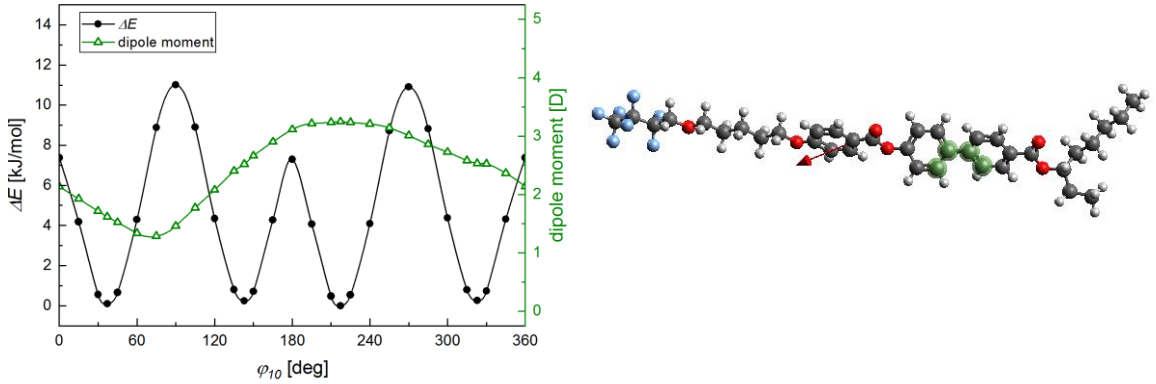


h)

Figure S3 (part II). Conformational energy (circles, left axis) and total dipole moment (squares, right axis) vs. torsional angle for selected intra-molecular rotations calculated for the isolated II.5.(HH) (S) molecule with the DFT-B3LYP/def2TZVP method.



i)



j)

Figure S3 (part III). Conformational energy (circles, left axis) and total dipole moment (squares, right axis) vs. torsional angle for selected intra-molecular rotations calculated for the isolated II.5.(HH) (S) molecule with the DFT-B3LYP/def2TZVP method.