## Probing kinetic and mechanistic features of bulk azide-alkyne cycloaddition

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Figure S1. <sup>1</sup>H NMR spectrum of 1-azidodecane (CDCl<sub>3</sub>-d1, 25 °C).



Figure S2. <sup>1</sup>H NMR spectrum of phenyl propargyl ether (CDCl<sub>3</sub>-d1, 25 °C).



**Figure S3**. <sup>1</sup>H NMR spectrum of 1-decyl-4-(phenoxymethyl)-1,2,3-triazole (1,4-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



**Figure S4**. <sup>13</sup>C NMR spectrum of compound 1-decyl-4-(phenoxymethyl)-1,2,3-triazole (1,4-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



**Figure S5.** NOESY <sup>1</sup>H-<sup>1</sup>H NMR spectrum of 1-decyl-4-(phenoxymethyl)-1,2,3-triazole (1,4-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



Figure S6. FTIR spectrum of 1-decyl-4-(phenoxymethyl)-1,2,3-triazole (1,4-adduct).



**Figure S7**. NMR <sup>1</sup>H spectrum of 1-decyl-5-(phenoxymethyl)-1,2,3-triazole (1,5-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



**Figure S8**. NMR <sup>13</sup>C spectrum of 1-decyl-5-(phenoxymethyl)-1,2,3-triazole (1,5-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



**Figure S9.** NOESY <sup>1</sup>H-<sup>1</sup>H NMR spectrum of 1-decyl-5-(phenoxymethyl)-1,2,3-triazole (1,5-adduct) (CDCl<sub>3</sub>-d1, 25 °C).



Figure S10. FTIR spectrum of 1-decyl-5-(phenoxymethyl)-1,2,3-triazole (1,5-adduct).



**Figure S11**. Original DSC curves for studied AAC (numbers denote heating rates in °C min-1) and *I*-azidodecane decomposition (B) (4 °C min $^{-1}$ ).



Figure S12. Chromatogram of the reaction mixture after DSC experiment at 0.5 °C min<sup>-1</sup>.



Figure S13. Chromatogram of the reaction mixture after DSC experiment at 1.0 °C min<sup>-1</sup>.



Figure S14. Chromatogram of the reaction mixture after DSC experiment at 2.0 °C min<sup>-1</sup>.



Figure S15. Chromatogram of the reaction mixture after DSC experiment at 4.0 °C min<sup>-1</sup>.



Figure S16. Calibration curve for the *1*,4-adduct.



Figure S17. Calibration curve for the *1*,5-adduct.