Supporting Information for

A reversible conductivity modulation of azobenzene-based ionic liquids in aqueous solutions by UV/Vis light

Zhiyong Li^{a,b}, Xiaoqing Yuan^{a,b}, Ying Feng^{a,b}, Yongkui Chen^{a,b}, Yuling Zhao^{a,b}, Huiyong Wang^{a,b}, Qingli Xu^{a,b}, Jianji Wang^{a,b*}

^a Collaborative Innovation Center of Henan Province for Green Manufacturing of Fine Chemicals, Key Laboratory of Green Chemical Media and Reactions, Ministry of Education, School of Chemistry and Chemical Engineering, Henan Normal University, Xinxiang, Henan 453007, P. R. China

^b Henan Key Laboratory of Green Chemistry, Henan Normal University, Xinxiang, Henan 453007, P. R. China

* E-mail: jwang@htu.cn

Content

- 1. ¹H NMR spectra of the azobenzene-based ionic liquids and related compounds
- 2. ¹H NMR data of the azobenzene-based ionic liquids and related compounds
- 3. UV-vis spectra for 0.01 mol·kg⁻¹ of the ionic liquids in aqueous solution

1. ¹H NMR spectra of the azobenzene-based ionic liquids and related compounds



Figure S1. ¹H NMR spectrum of C₄Azo



Figure S2. ¹H NMR spectrum of C_4AzoC_2Br



Figure S3. ¹H NMR spectrum of C_4AzoC_4Br







Figure S5. ¹H NMR spectrum of $[C_4AzoC_2DMEA]Br$



Figure S6. ¹H NMR spectrum of [C₄AzoC₄DMEA]Br



Figure S7. ¹H NMR spectrum of $[C_4AzoC_6DMEA]Br$



Figure S8. ¹H NMR spectrum of [C₄AzoC₂TMA]Br



Figure S9. ¹H NMR spectrum of [C₄AzoC₂MIM]Br

2. ¹H NMR data of the azobenzene-based ionic liquids and related compounds

1. **C₄AzoOH:** ¹H NMR (400MHz, DMSO-d₆, TMS): δ=10.28(s, 1H, -OH), 7.79(d, 2H, Ph-H), 7.73(d, 2H, Ph-H), 7.38(d, 2H, Ph-H), 6.95(d, 2H, Ph-H), 2.66 (t, 2H, -CH₂), 1.59 (m, 2H, -CH₂), 1.32 (m, 2H, -CH₂), 0.91 (t, 3H, -CH₃) ppm

2. C₄AzoC₂Br: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.87(d, 2H, Ph-H), 7.78(d, 2H, Ph-H), 7.38(d, 2H, Ph-H), 7.15(d, 2H, Ph-H), 4.44 (t, 2H, -CH₂), 3.85 (t, 2H, -CH₂), 2.68 (t, 2H, -CH₂), 1.60 (m, 2H, -CH₂), 1.33 (m, 2H, -CH₂), 0.91 (t, 3H, -CH₃) ppm

3. C₄AzoC₄Br: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.85(d, 2H, Ph-H), 7.76(d, 2H, Ph-H), 7.37(d, 2H, Ph-H), 7.13(d, 2H, Ph-H), 4.11 (t, 2H, -CH₂), 3.62 (t, 2H, -CH₂), 2.65 (t, 2H, -CH₂), 2.01 (m, 2H, -CH₂), 1.87 (m, 2H, -CH₂), 1.61 (m, 2H, -CH₂), 1.34 (m, 2H, -CH₂), 0.91 (t, 3H, -CH₃) ppm

4. **C₄AzoC₆Br**: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.87(d, 2H, Ph-H), 7.75(d, 2H, Ph-H), 7.40(d, 2H, Ph-H), 7.12(d, 2H, Ph-H), 4.08 (t, 2H, -CH₂), 3.55 (t, 2H, -CH₂), 2.65 (t, 2H, -CH₂), 1.82 (m, 2H, -CH₂), 1.75 (m, 2H, -CH₂), 1.60 (m, 2H, -CH₂), 1.46 (m, 4H, -CH₂), 1.34 (m, 2H, -CH₂), 0.91 (t, 3H, -CH₃) ppm

5. **[C₄AzoC₂DMEA]Br**: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.92(d, 2H, Ph-H), 7.79(d, 2H, Ph-H), 7.41(d, 2H, Ph-H), 7.21(d, 2H, Ph-H), 5.34 (t, 1H, -OH), 4.59 (t, 2H, -CH₂), 3.90

(t, 4H, -CH₂), 3.57 (t, 2H, -CH₂), 3.20 (s, 6H, -CH₃), 2.69 (t, 2H, -CH₂), 1.60 (m, 2H, -CH₂), 1.34 (m, 2H, -CH₂), 0.93 (t, 3H, -CH₃) ppm

6. **[C₄AzoC₄DMEA]Br**: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.87(d, 2H, Ph-H), 7.76(d, 2H, Ph-H), 7.41(d, 2H, Ph-H), 7.16(d, 2H, Ph-H), 5.31 (t, 1H, -OH), 4.16 (t, 2H, -CH₂), 3.84 (t, 2H, -CH₂), 3.41 (t, 4H, -CH₂), 3.09 (s, 6H, -CH₃), 2.67 (t, 2H, -CH₂), 1.87 (m, 2H, -CH₂), 1.80 (m, 2H, -CH₂), 1.64 (m, 2H, -CH₂), 1.34 (m, 2H, -CH₂), 0.93 (t, 3H, -CH₃) ppm

[C₄AzoC₆DMEA]Br: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.85(d, 2H, benzene ring-H), 7.75(d, 2H, benzene ring-H), 7.39(d, 2H, benzene ring-H), 7.13(d, 2H, benzene ring-H), 5.30 (t, 1H, -OH), 4.08 (t, 2H, -CH₂), 3.82 (m, 2H, -CH₂), 3.40 (m, 2H, -CH₂), 3.32 (m, 2H, -CH₂), 3.07 (s, 6H, -NCH₃), 2.66 (t, 2H, -CH₂), 1.78 (m, 4H, -CH₂), 1.59 (m, 2H, -CH₂), 1.47 (m, 2H, -CH₂), 1.31 (m, 4H, -CH₂), 0.90 (t, 3H, -CH₃) ppm

8. [C₄AzoC₂TMA]Br: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=7.91(d, 2H, Ph-H), 7.78(d, 2H, Ph-H), 7.41(d, 2H, Ph-H), 7.22(d, 2H, Ph-H), 4.59 (t, 2H, -CH₂), 3.85 (t, 2H, -CH₂), 3.20 (s, 9H, -CH₃), 2.67 (m, 2H, -CH₂), 1.60 (m, 2H, -CH₂), 1.33 (m, 2H, -CH₂), 0.91 (t, 3H, -CH₃) ppm

9. **[C₄AzoC₂MIM]Br**: ¹H NMR (400MHz, DMSO-d₆, TMS): δ=9.26(s, 1H, imidazole ring-H), 7.89(t, 2H, Ph-H, and 1H, imidazole ring-H), 7.78(t, 2H, Ph-H, and 1H, imidazole ring-H), 7.40(d, 2H, Ph-H), 7.18(d, 2H, Ph-H), 4.67 (t, 2H, -CH₂), 4.49 (t, 2H, -CH₂), 3.90(s, 3H,-NCH₃), 2.67 (t, 2H, -CH₂), 1.61 (m, 2H, -CH₂), 1.34 (m, 2H, -CH₂), 0.93 (t, 3H, -CH₃) ppm

3. UV-vis spectra for 0.01 mol·kg⁻¹ of the ionic liquids in aqueous solution



Figure S10. UV-vis spectra for aqueous solution of 0.01 mol·kg⁻¹ [C₄AzoC₂TMA]Br after different UV irradiation time at 25.0°C: a, initial state; b, 2min; c, 4 min; d, 6 min; e, 8 min; f, 10 min; g, 15 min; h, 20 min; i, 40min; j, 60min.



Figure S11. UV-vis spectra for aqueous solution of 0.01 mol·kg⁻¹ [C₄AzoC₂MIM]Br after different UV irradiation time at 25.0°C: a, initial state; b, 2min; c, 4 min; d, 6 min; e, 8 min; f, 10 min; g, 15 min; h, 20 min; i, 40min; j, 60min.



Figure S12. UV-vis spectra for aqueous solution of 0.01 mol·kg⁻¹ [C₄AzoC₄DMEA]Br after different UV irradiation time at 25.0°C: a, initial state; b, 2min; c, 4 min; d, 6 min; e, 8 min; f, 10 min; g, 15 min; h, 20 min; i, 40min; j, 60min.



Figure S13. UV-vis spectra for aqueous solution of 0.01 mol·kg⁻¹ [C₄AzoC₆DMEA]Br after different UV irradiation time at 25.0°C: a, initial state; b, 2min; c, 4 min; d, 6 min; e, 8 min; f, 10 min; g, 15 min; h, 20 min; i, 40min; j, 60min.