1 Supporting information

2 Ionic liquid-doped and p-NIPAAm-based copolymer (p-NIBIm): extraordinary drug-

3 entrapping and -releasing behaviors at 38-42 °C

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11 Experimental data

- 12 Fig. S1 ¹H-NMR spectra of [BVIm]Br, *p*-NIBIm, *p*-BVIm, and *p*-NIPAAm.
- 13 a) ¹H-NMR spectrum of [BVIm]Br showed characteristic peaks as follows:



¹H-NMR (*d*₆-DMSO): (ppm) 0.91 (t, 3H-9), 1.30 (m, 2H-8), 1.82 (m, 2H-7), 4.22 (q, 2H-6),
5.43~5.44 (dd, 1H-11), 5.95~5.97 (dd, 1H-11), 7.32 (dd, 1H-10), 7.96 (s, 1H-4), 8.22 (s, 1H-17),
5), 9.59 (s, 1H-2).



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3 b) ¹H-NMR spectrum of the *p*-NIBIm copolymer also showed characteristic peaks as follows:



¹H-NMR (*d*6-DMSO): (ppm) 0.90 (bt, 3H-9), 0.97~1.17 (bd, 6H-e), 1.28~1.39 (bm + bd, 2H-6 8 + 2H-11), 1.40~1.71 (bd, 2H-a), 1.73~1.88 (bm + bt, 2H-7 + 1H-10), 1.89~2.45 (bt, 1H-b),
3.84 (bm, 1H-d), 4.15 (bt, 2H-6), 6.90~7.64 (bt, 1H-c or 1NH), 7.65~7.95 (bd + bd, 1H-4 + 8 1H-5), 9.80 (bs, 1H-2).

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2 c) ¹H-NMR spectrum of the *p*-BVIm showed characteristic peaks as follows:



4 ¹H-NMR (*d*₆-DMSO): (ppm) 0.89 (bt, 3H-9), 1.27 (bm + bd, 2H-8 + 2H-11), 1.79 (bm + bt,

5 2H-7 + 1H-10), 4.1-4.3 (bt, 2H-6), 7.60~8.25 (bd + bd, 1H-4 + 1H-5), 9.4~9.8 (bs, 1H-2).



3 d) ¹H-NMR spectrum of the *p*-NIPAAm polymer showed characteristic peaks as follows:



- 5 ¹H-NMR (*d*₆-DMSO): (ppm) 1.05 (m, 6H-e), 1.20~1.71 (bd, 2H-a), 1.96 (bt, 1H-b), 3.85 (m,
- 6 1H-d), 6.95~7.65 (bt, 1H-c or NH).



- 3 Fig. S2 MALDI-TOF spectra of *p*-NIBIm, *p*-BVIm, and *p*-NIPAAm.
- 4 a) *p*-NIBIm
- 5 Mn = 1353; Mw = 2001; Mw/Mn = 1.47



8 b) *p*-BVIm





2 c) *p*-NIPAAm

3 Mn = 1998; Mw = 2521; Mw/Mn = 1.26



6 Fig. S3 FT-IR spectra of *p*-NIBIm, *p*-BVIm, and *p*-NIPAAm.





Fig. S4 LCST determination of IL-doped *p*-NIBIm using DSC scan; a) via several continuous
 heating and cooling cycles at pH=7 and b) in the acidic pH environments (pH=4-5).

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5 Fig. S5 Colloidal stability of the aqueous *p*-NIBIm solution (0.5 mg/mL) at 37 °C.



1 Fig. S6 Temperature-dependant size change of the *p*-NIPAAm polymer and the complex with



2 BSA, *p*-NIPAAm/BSA.

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- 4
- 5 Fig. S7 The calibration curve, which is created by plotting the known BSA concentration on x
- 6 axis and the absorbance of BSA/DC complexes at λ_{max} =750 nm on y axis.



BSA calibration-curve