

Supplementary Materials:

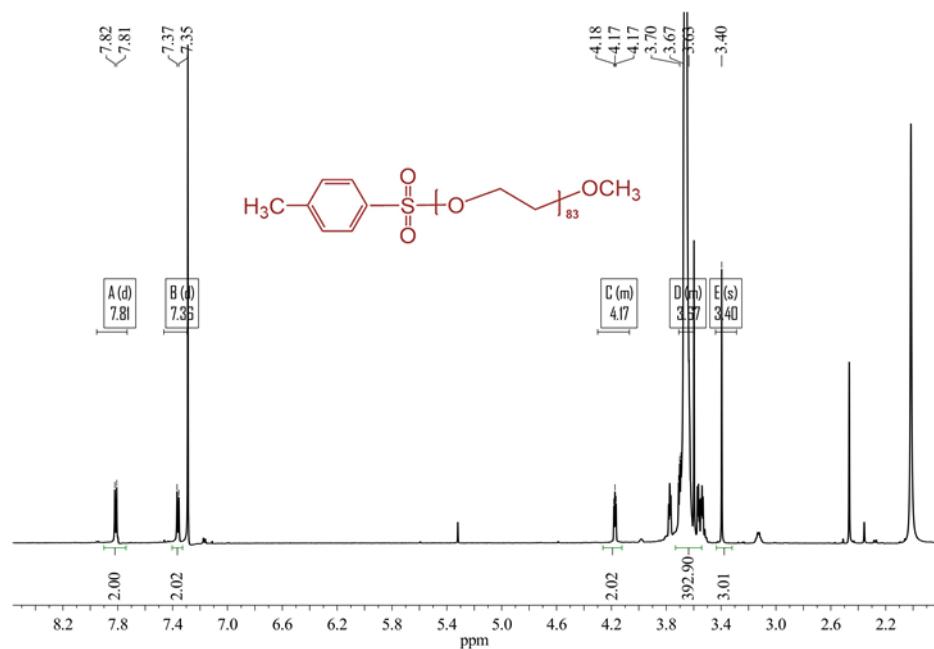


Figure S1. ¹H-NMR spectrum of PEG4k-TsCl.

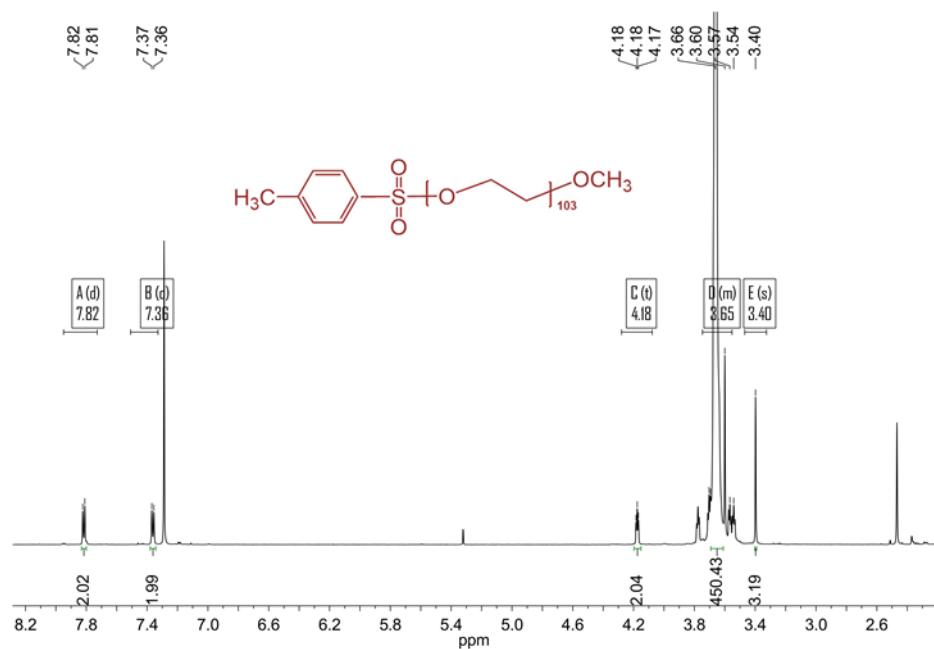


Figure S2. ¹H-NMR spectrum of PEG5k-TsCl.

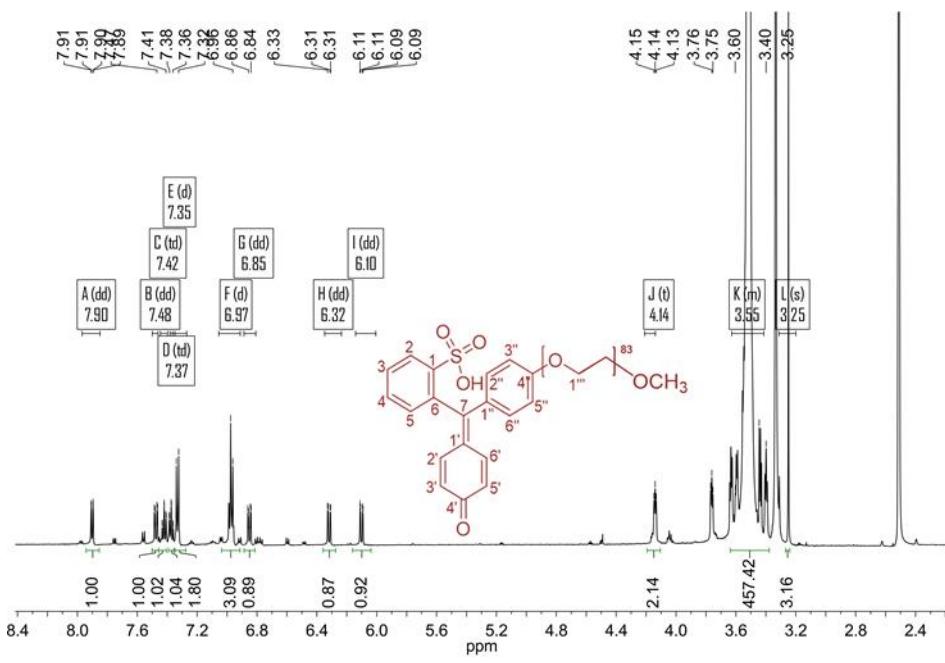


Figure S3. ¹H-NMR spectrum of PEG4k-PR.

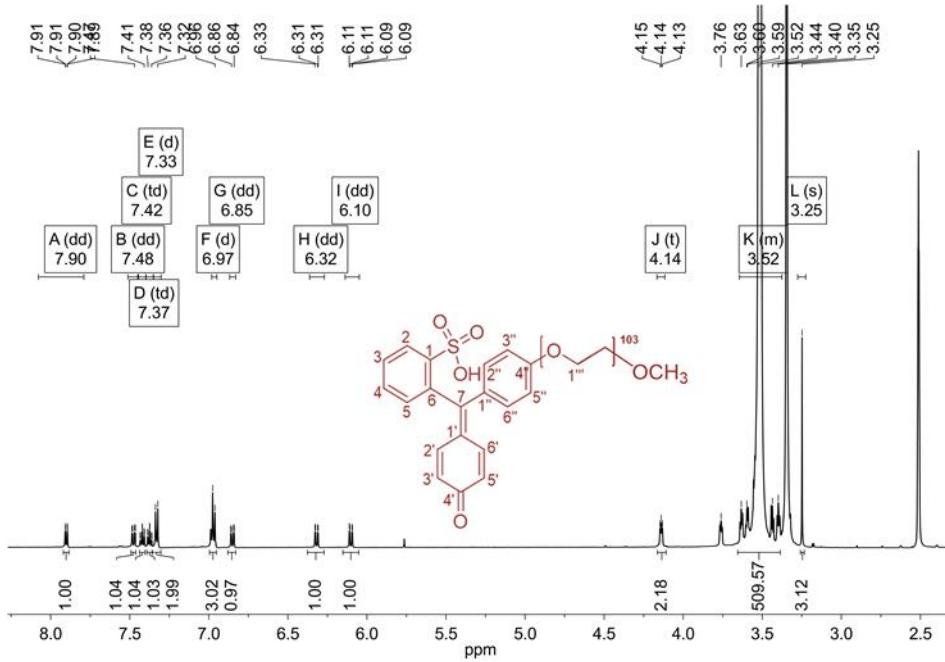


Figure S4. ¹H-NMR spectrum of PEG5k-PR.

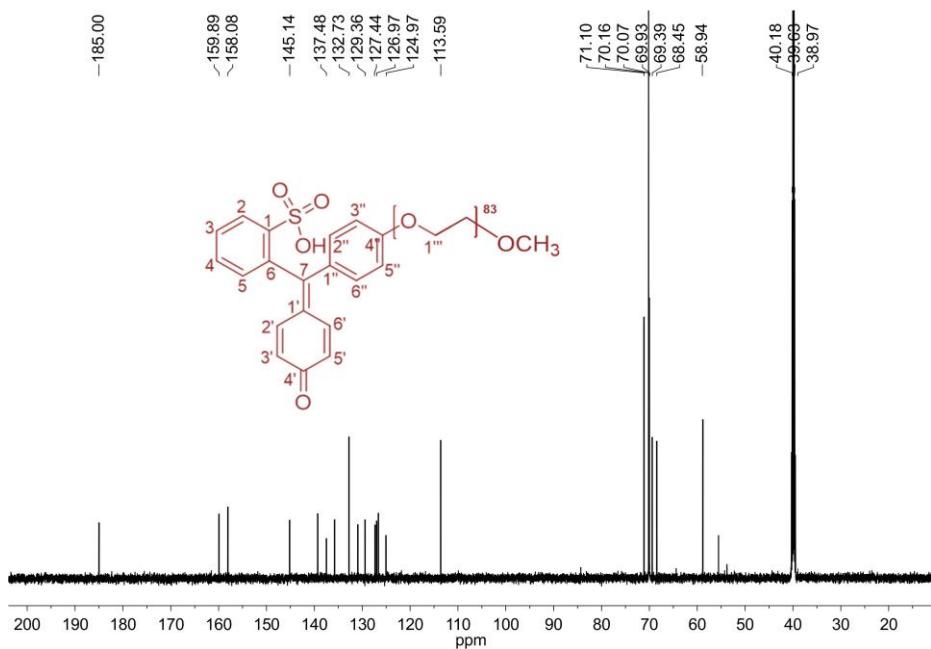


Figure S5. ^{13}C -NMR spectrum of PEG4k-PR.

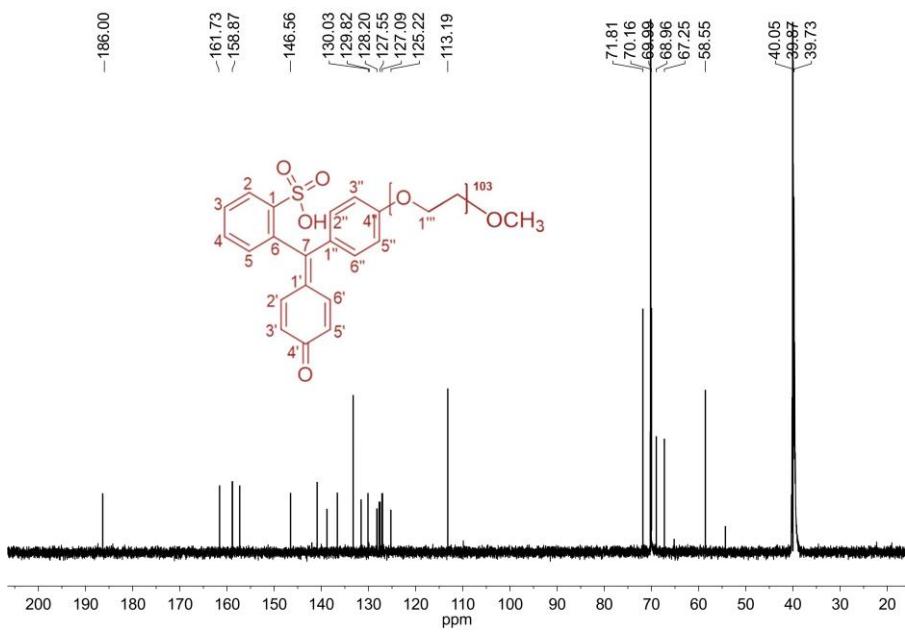


Figure S6. The ^{13}C -NMR spectrum of PEG5k-PR.

Table S1. Concentration changes of all compounds before and after the shaker bath experiment. Data are shown as mean \pm SD, n = 3.

Compound	Initial		Ending (120 min)		Percentage of concentration change
	As/Ai	Conc. (μ M)	As/Ai	Conc. (μ M)	
PR	0.32 \pm 0.01	56.90 \pm 1.10	0.31 \pm 0.01	56.11 \pm 1.57	1.38%
PEG4kPR	0.54 \pm 0.01	54.87 \pm 2.44	0.54 \pm 0.01	54.29 \pm 0.61	1.07%
PEG5k-PR	0.87 \pm 0.01	56.47 \pm 0.65	0.86 \pm 0.01	55.58 \pm 0.37	1.57%
Atenolol	0.71 \pm 0.01	105.19 \pm 0.96	0.71 \pm 0.01	105.93 \pm 0.61	0.75
Ketoprofen	1.11 \pm 0.02	101.27 \pm 1.22	1.09 \pm 0.01	100.28 \pm 0.96	0.97
Metoprolol	0.79 \pm 0.02	103.62 \pm 2.29	0.76 \pm 0.01	102.12 \pm 0.58	1.46

As: HPLC peak area of compound.

Ai: HPLC peak area of internal standard.

Percentage of concentration change: (Initial Conc.-Ending Conc.) / Initial Conc.

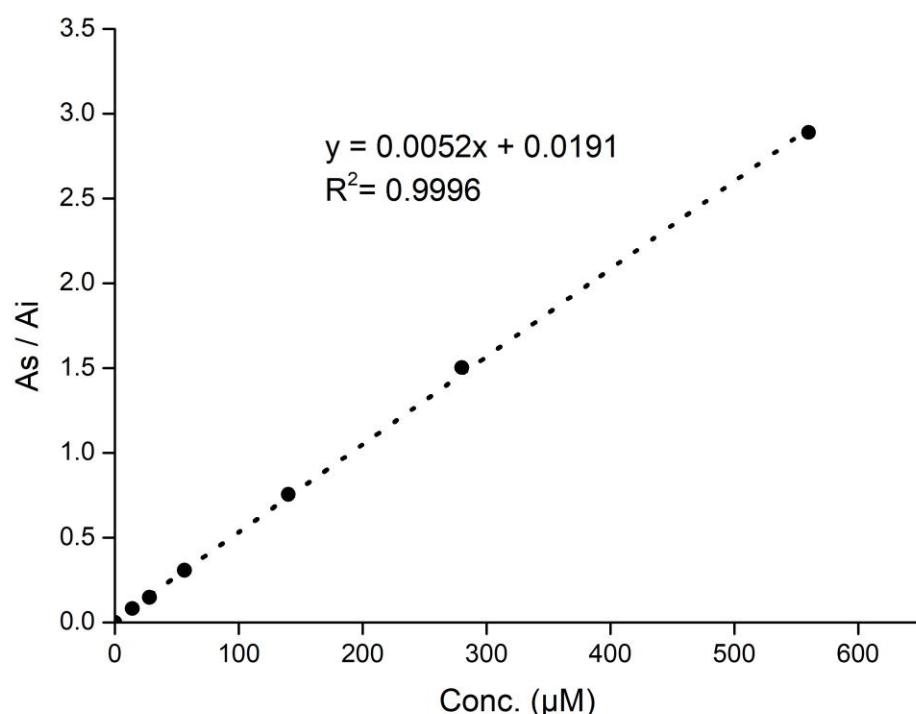


Figure S7. Calibration line of PR (MW = 354) in Krebs–Ringer buffer.

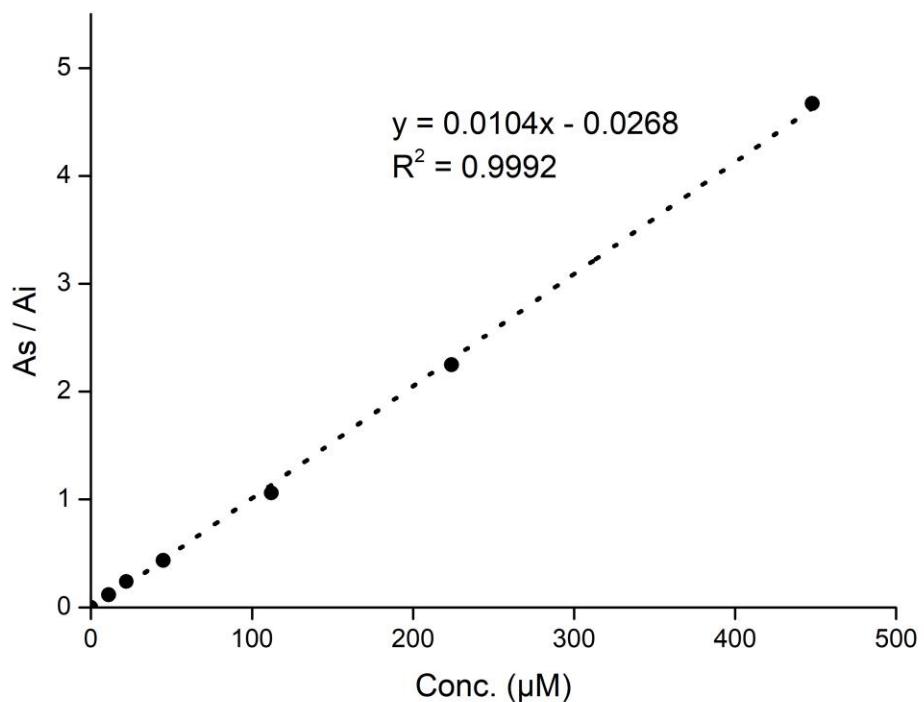


Figure S8. Calibration line of PEG4k-PR (MW = 4354) in Krebs–Ringer buffer.

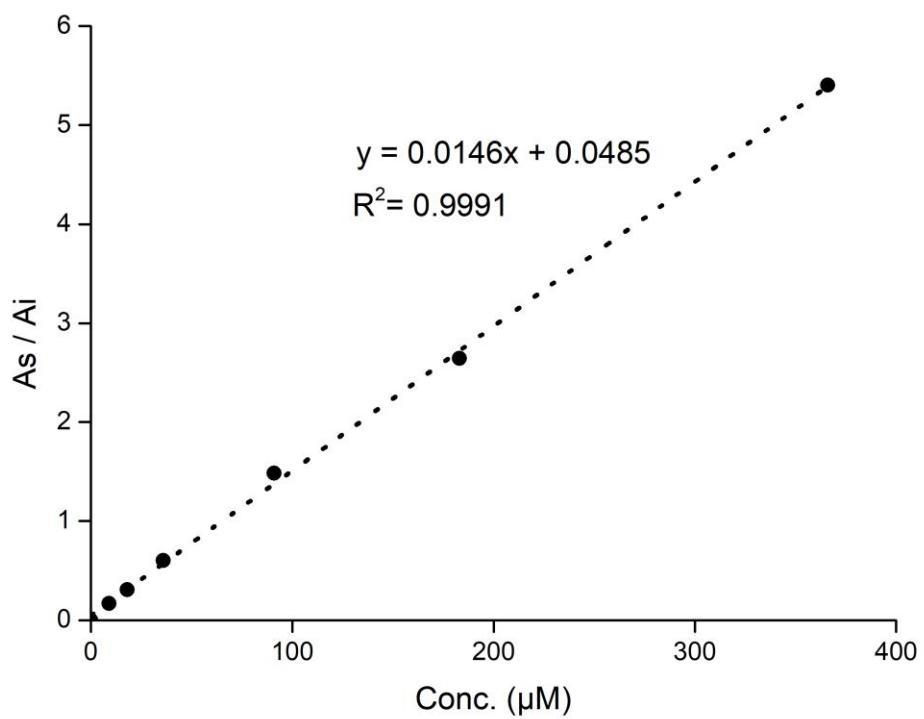


Figure S9. Calibration line of PEG5k-PR (MW = 5354) in Krebs–Ringer buffer.

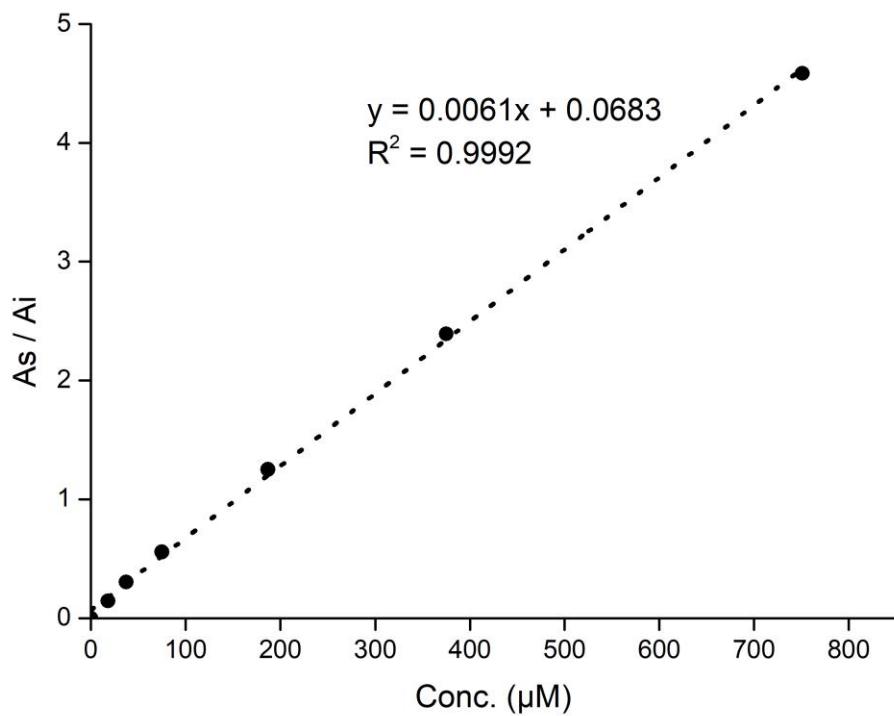


Figure S10. Calibration line of atenolol (MW = 266) in Krebs–Ringer buffer.

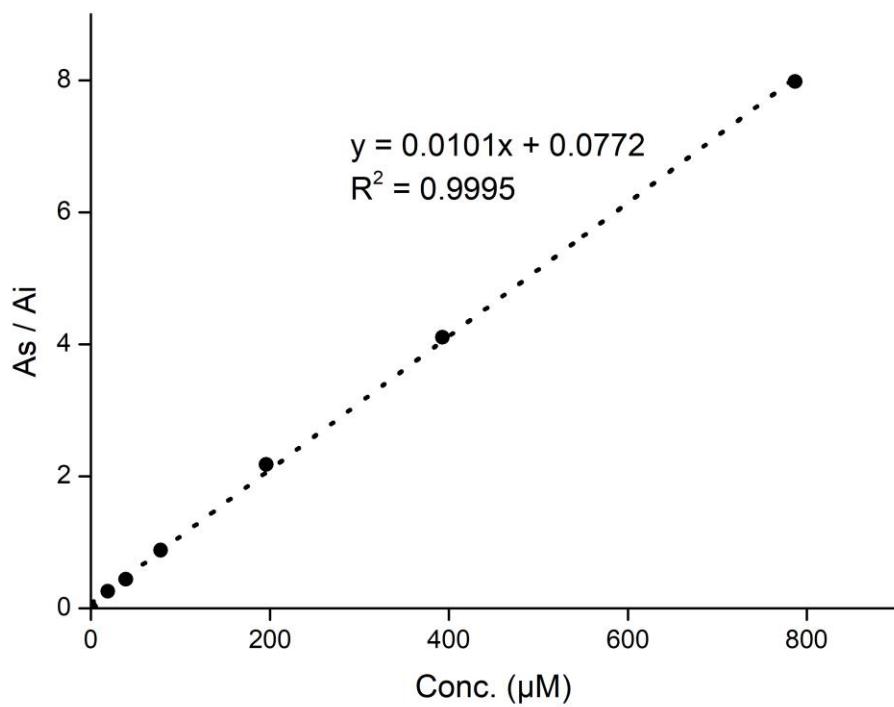


Figure S11. Calibration line of ketoprofen (MW = 254) in Krebs–Ringer buffer.

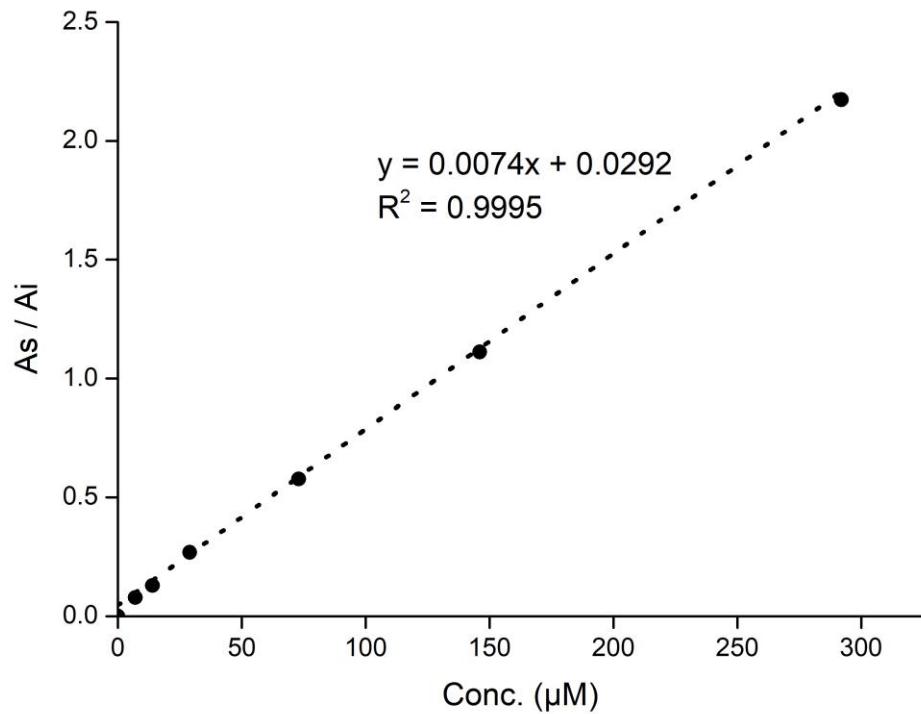


Figure S12. Calibration line of metoprolol (MW = 684) in Krebs–Ringer buffer.