

Supplementary File

NMR analysis to confirm π - π interactions, aromatic hydrophobic interactions:

Commercial samples of poly(NaSS), \bar{M} ~70,000 Da obtained from Sigma-Aldrich and poly(DMA), \bar{M} ~100,000 Da, obtained from Scientific Polymer Products, Inc. (Ontario, NY) were used for proton NMR analysis. Aromatic protons of MB are labelled as H₁, H₂ and H₃, which showed broadening of signals in the presence of aromatic polymer like poly(NaSS) whereas no peak broadening of signals observed in poly(DMA) solutions. Signal broadening confirmed the π - π interaction and delocalization of electron clouds as part of aromatic hydrophobic interactions between poly(NaSS) and MB. Signal broadening of MB proton signals was not observed for co-solutions of poly(DMA) and MB.

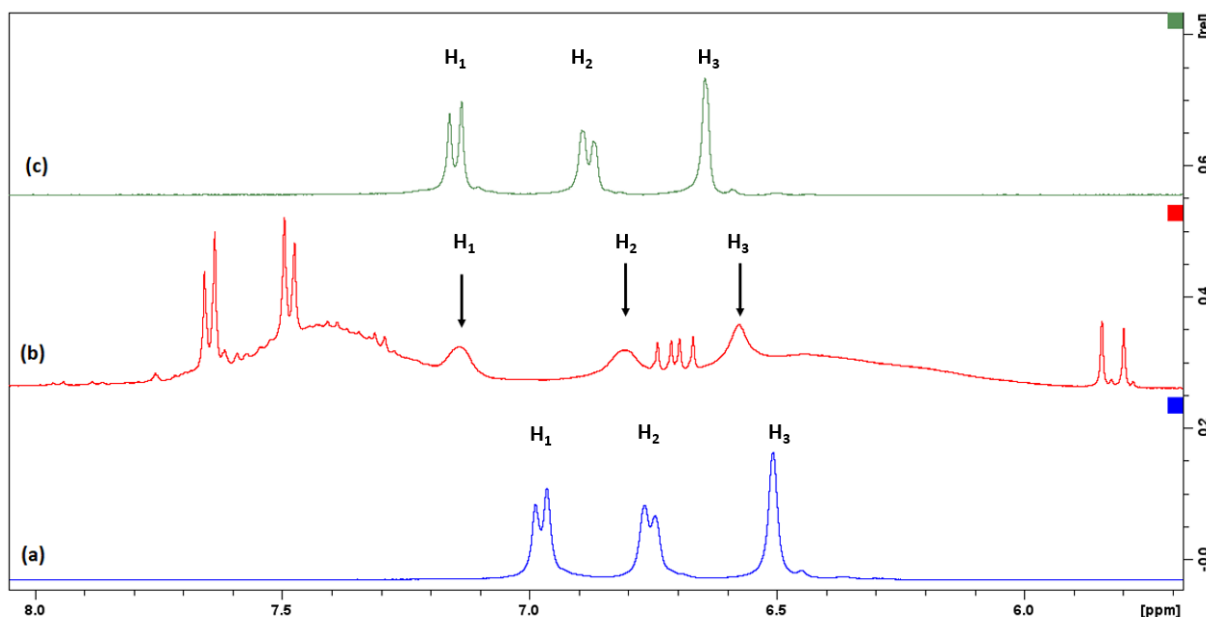


Figure S1- ¹H NMR spectra comparison for (a) 10⁻³ M of MB solution in D₂O, (b) 10⁻³ M of MB in solution with 10⁻² M poly (NaSS) and (c) 10⁻³ M of MB in solution with 10⁻² M poly (DMA)

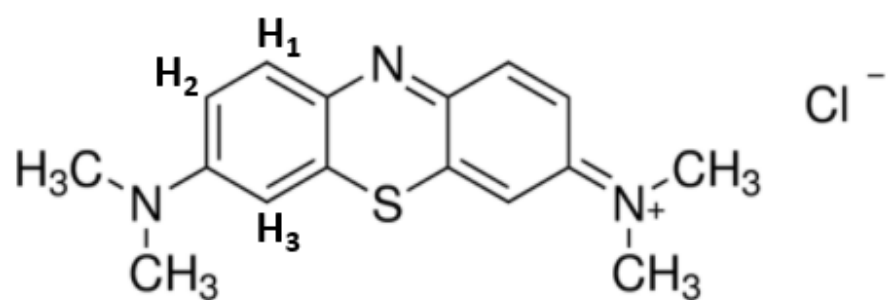


Figure S2 – Chemical structure of MB dye