

Supplementary Information

(Co)Polymers Containing Boron Difluoride 3-
Cyanofornazanate Complexes: Emission Enhancement via
Random Copolymerization

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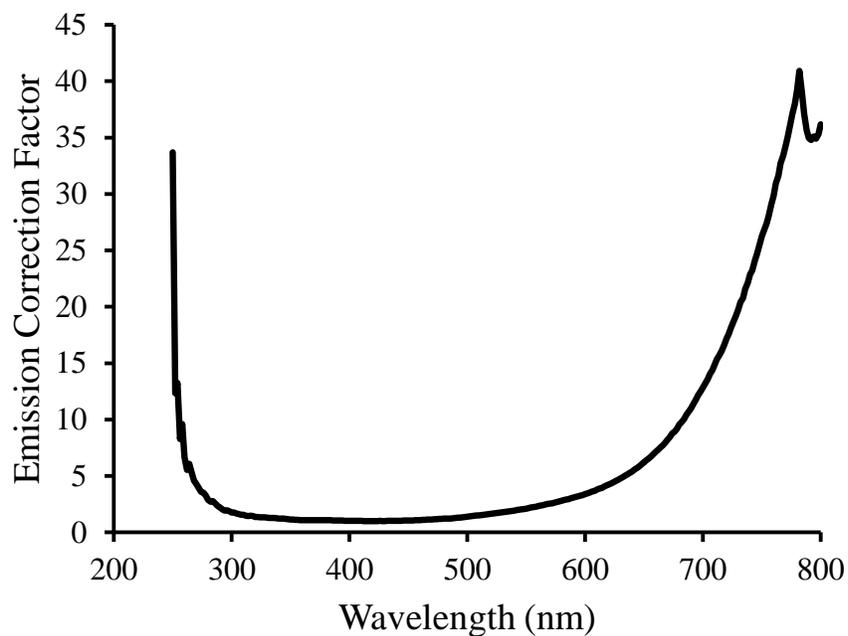


Fig. S1 Wavelength-dependent emission correction provided by Photon Technology International.

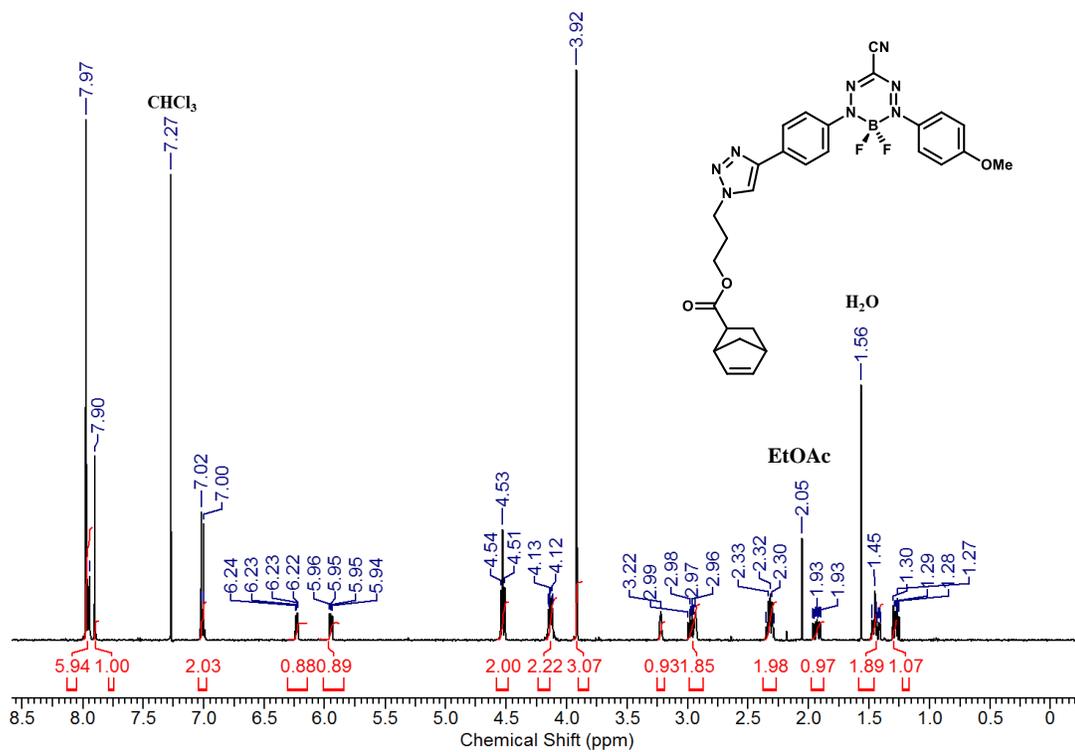


Fig. S2 ¹H NMR spectrum of monomer **BF2N** in CDCl₃.

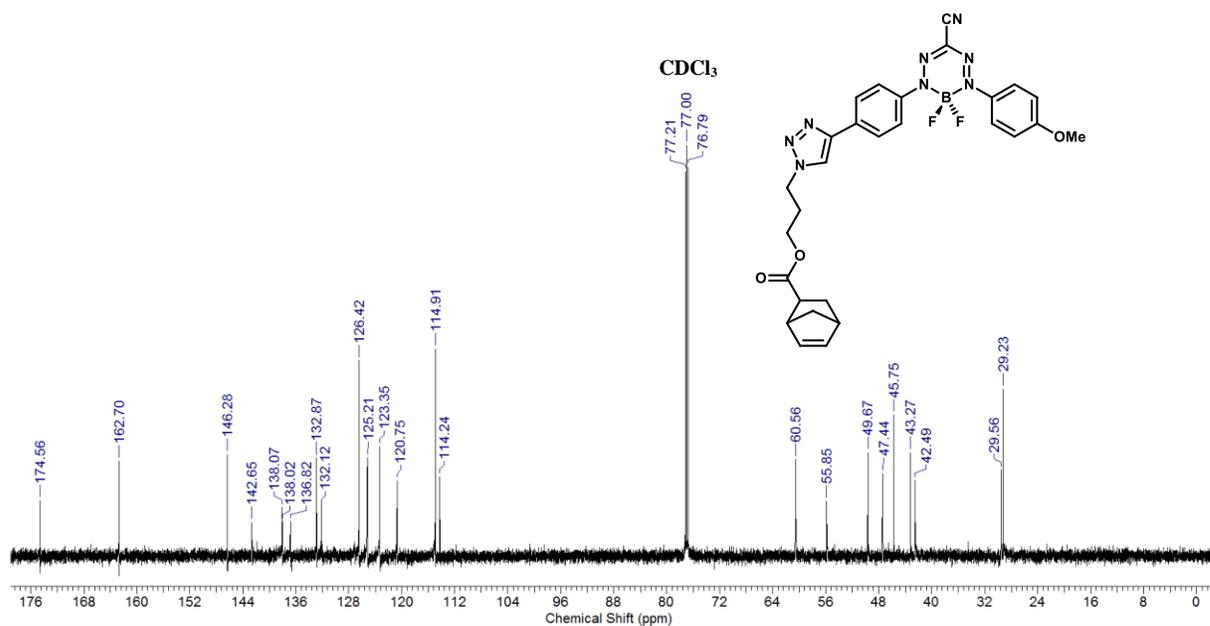


Fig. S3 $^{13}\text{C}\{^1\text{H}\}$ NMR spectrum of monomer **BF2N** in CDCl_3 .

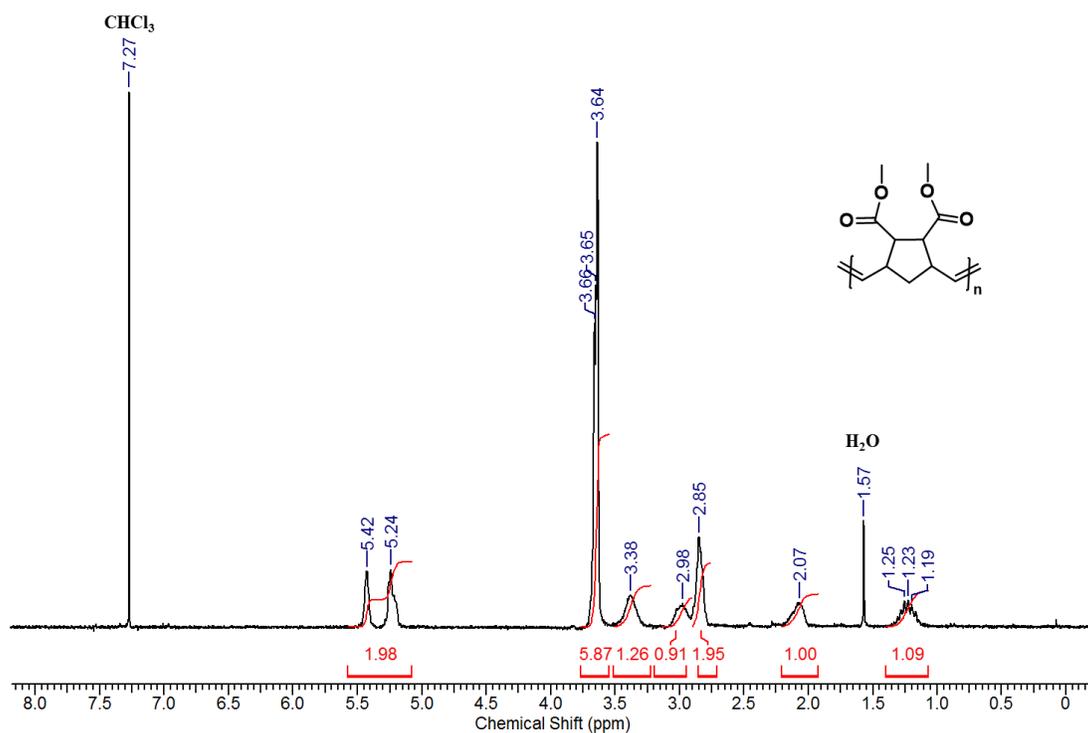


Fig. S4 ^1H NMR spectrum of **PDND** in CDCl_3 .

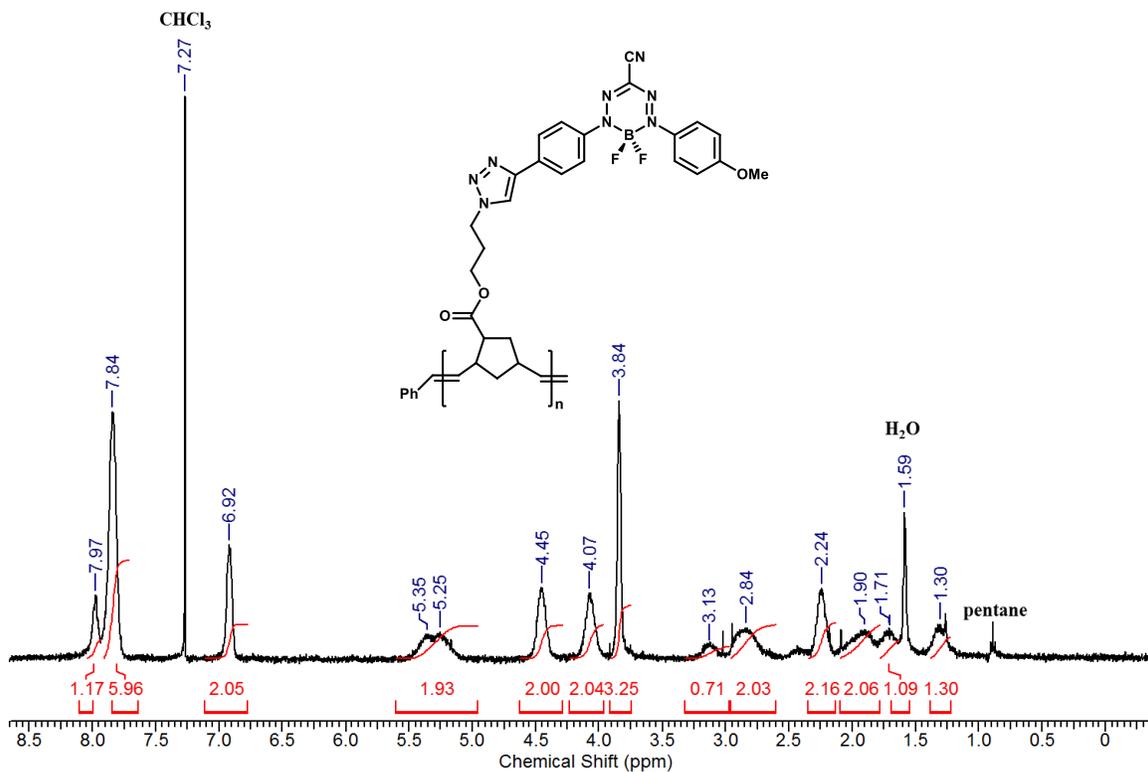


Fig. S5 ^1H NMR spectrum of PBF2N in CDCl_3 .

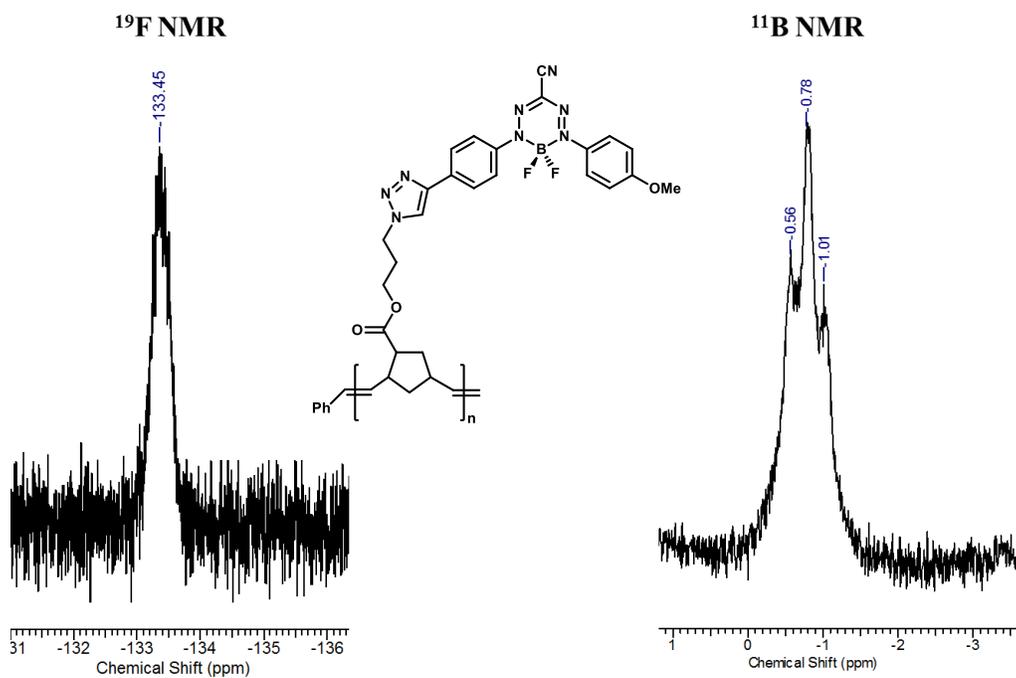


Fig. S6 ^{19}F NMR (left) spectrum and ^{11}B NMR (right) of PBF2N in CDCl_3 .

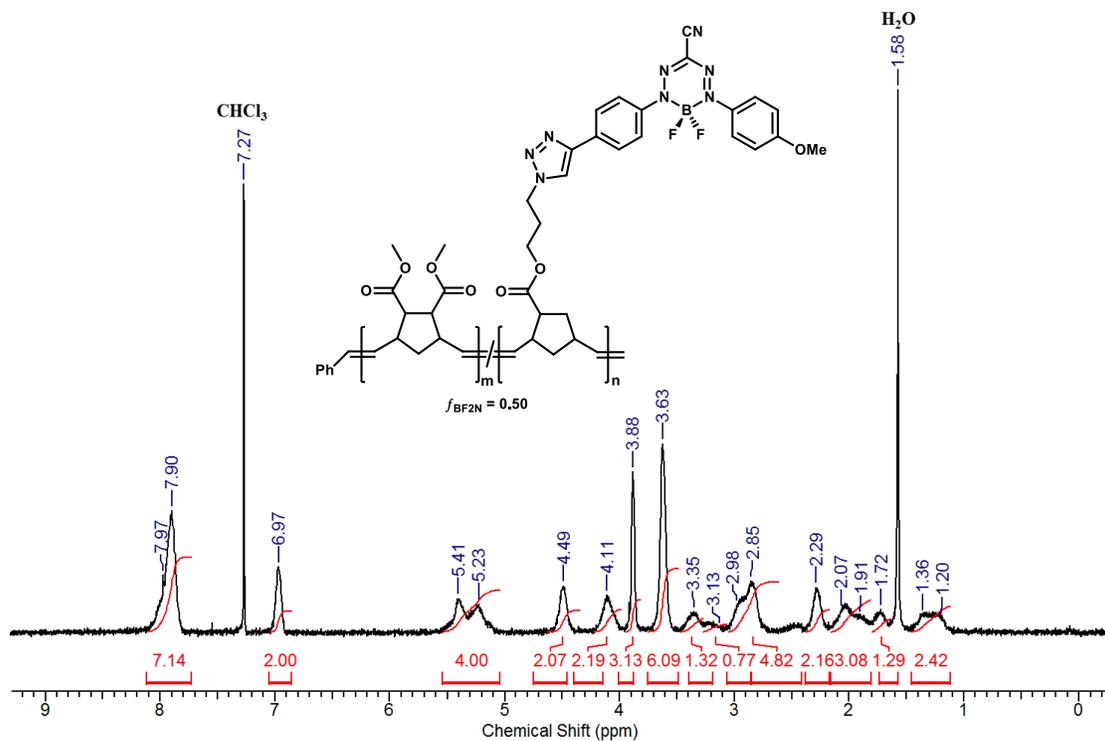


Fig. S7 1H NMR spectrum of $(PDND)_m-r-(PBF_2N)_n$ ($f_{BF_2N} = 0.50$) in $CDCl_3$.

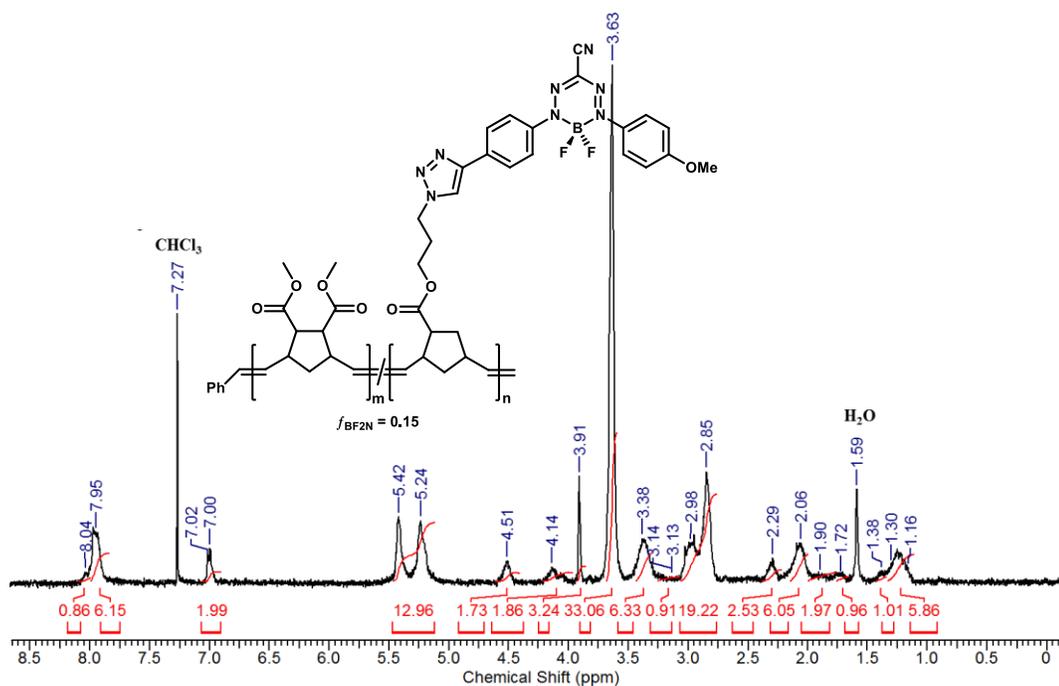


Fig. S8 1H NMR spectrum of $(PDND)_m-r-(PBF_2N)_n$ ($f_{BF_2N} = 0.15$) in $CDCl_3$.

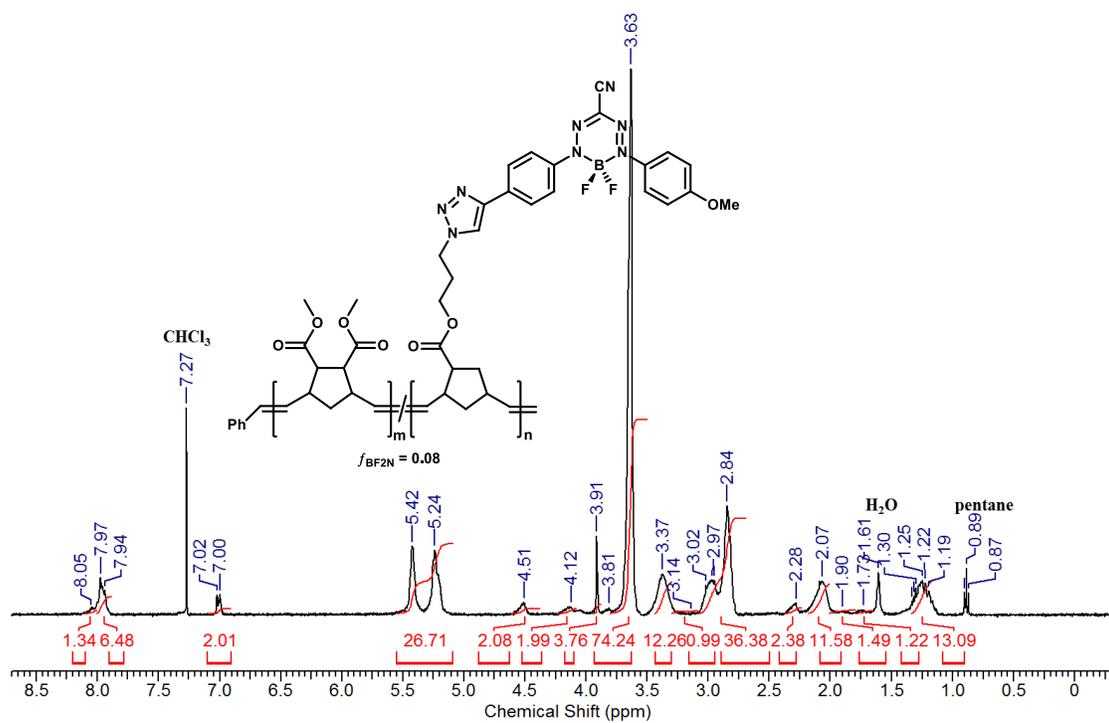


Fig. S9 ^1H NMR spectrum of $(\text{PDND})_m\text{-}r\text{-}(\text{PBF}_2\text{N})_n$ ($f_{\text{BF}_2\text{N}} = 0.08$) in CDCl_3 .

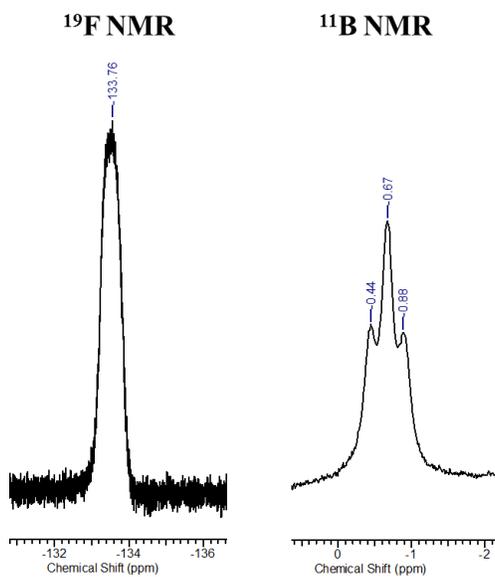


Fig. S10 Representative ^{11}B NMR and ^{19}F NMR for random copolymers $(\text{PDND})_m\text{-}r\text{-}(\text{PBF}_2\text{N})_n$ and block copolymers $(\text{PDND})_m\text{-}b\text{-}(\text{PBF}_2\text{N})_n$ in CDCl_3 .

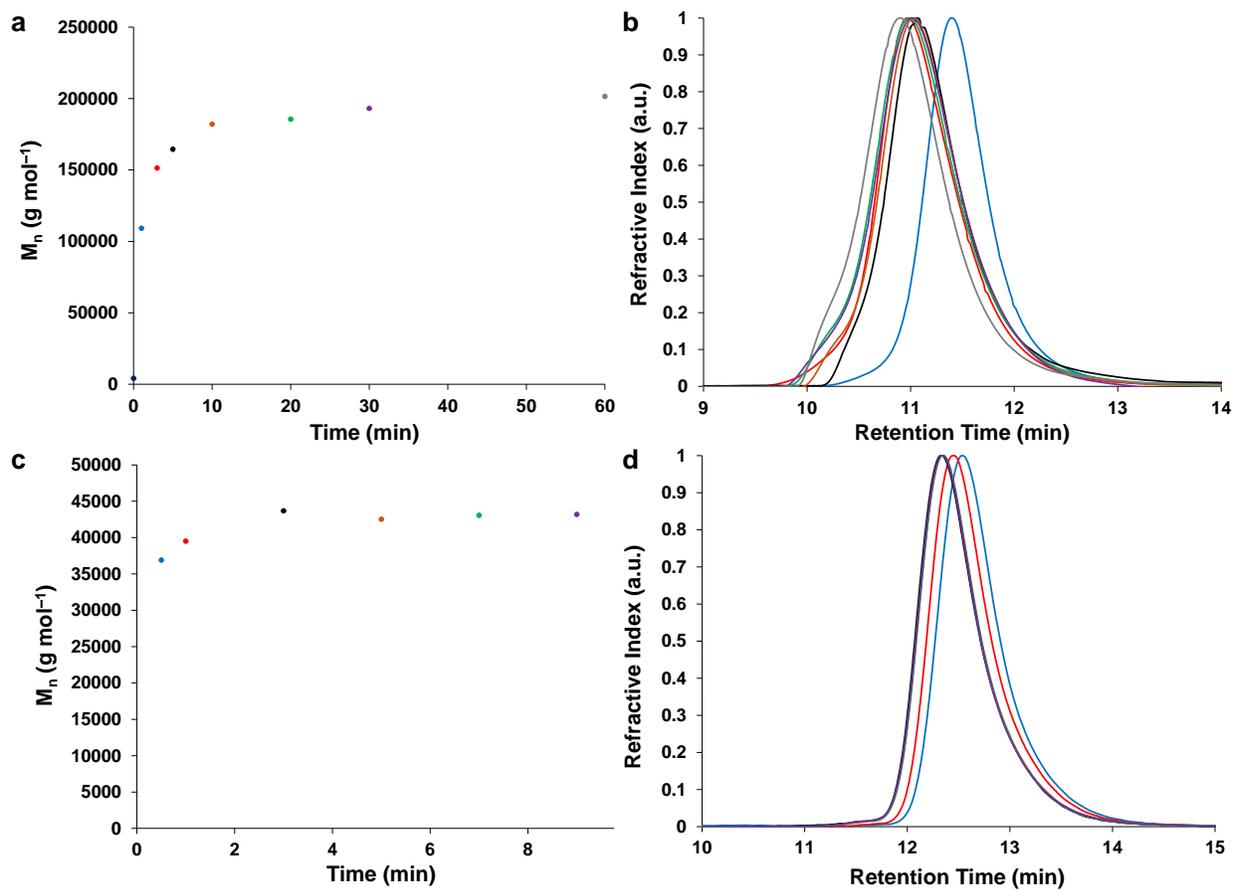


Fig. S11 Relationship between number average molecular weight (M_n) of homopolymers **PBF2N** (a,b) and **PDND** (c,d) and reaction time.

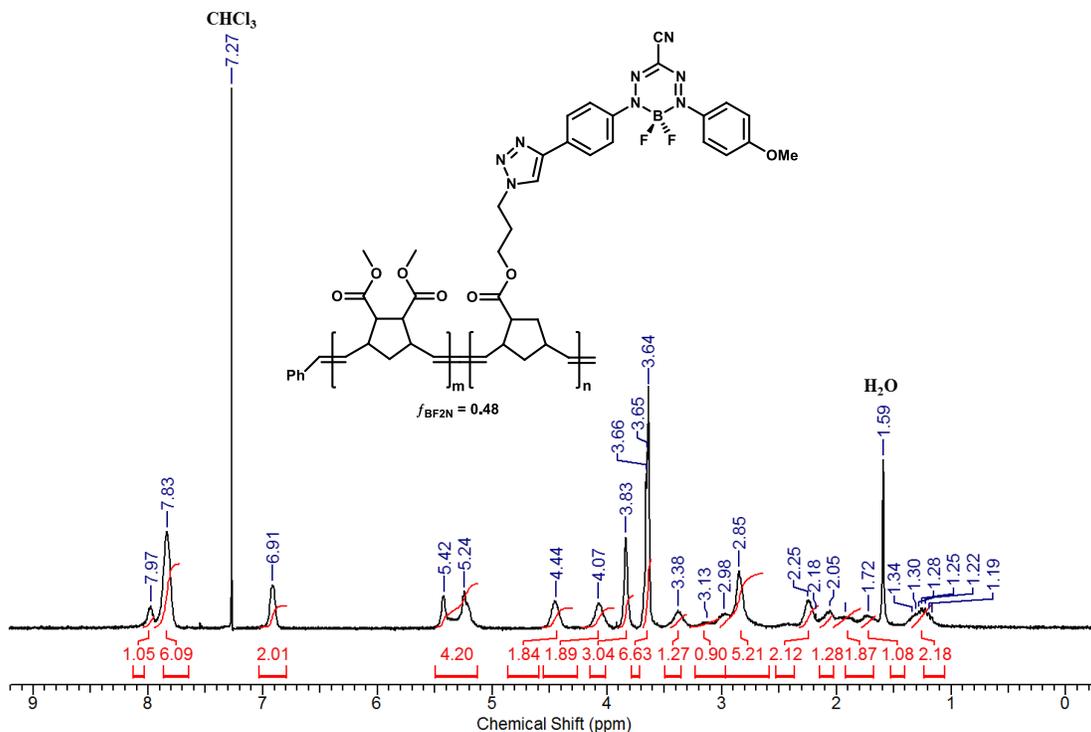


Fig. S12 $^1\text{H NMR}$ spectrum of $(\text{PDND})_m\text{-}b\text{-(PBF}_2\text{N)}_n$ ($f_{\text{BF}_2\text{N}} = 0.48$) in CDCl_3 .

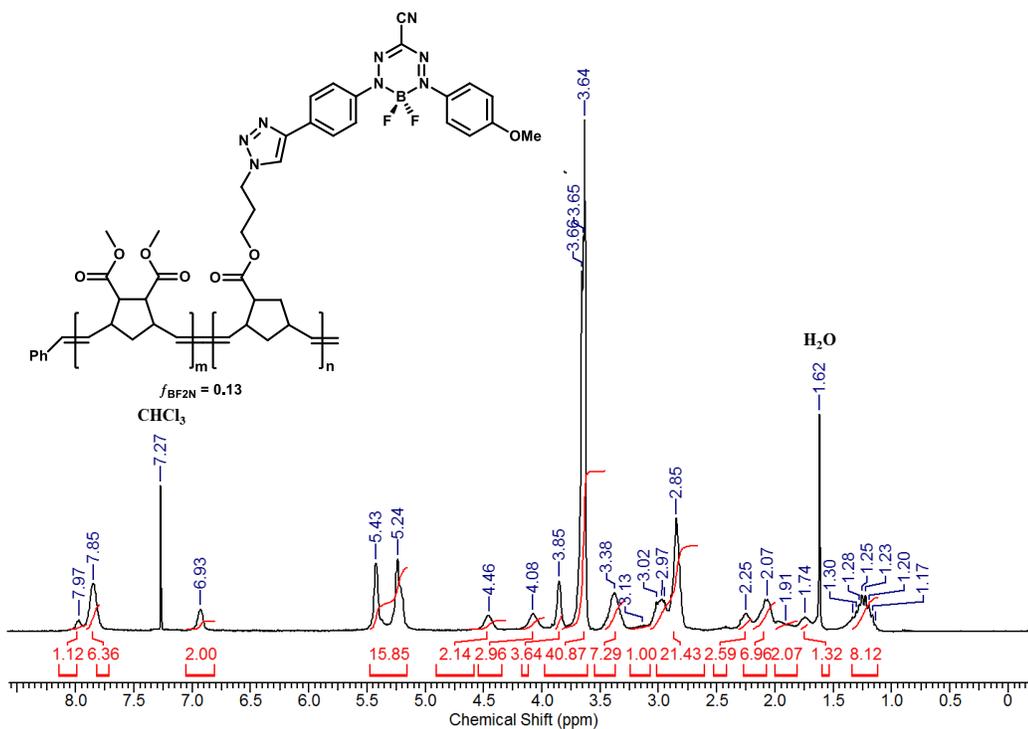


Fig. S13 $^1\text{H NMR}$ spectrum of $(\text{PDND})_m\text{-}b\text{-(PBF}_2\text{N)}_n$ ($f_{\text{BF}_2\text{N}} = 0.13$) in CDCl_3 .

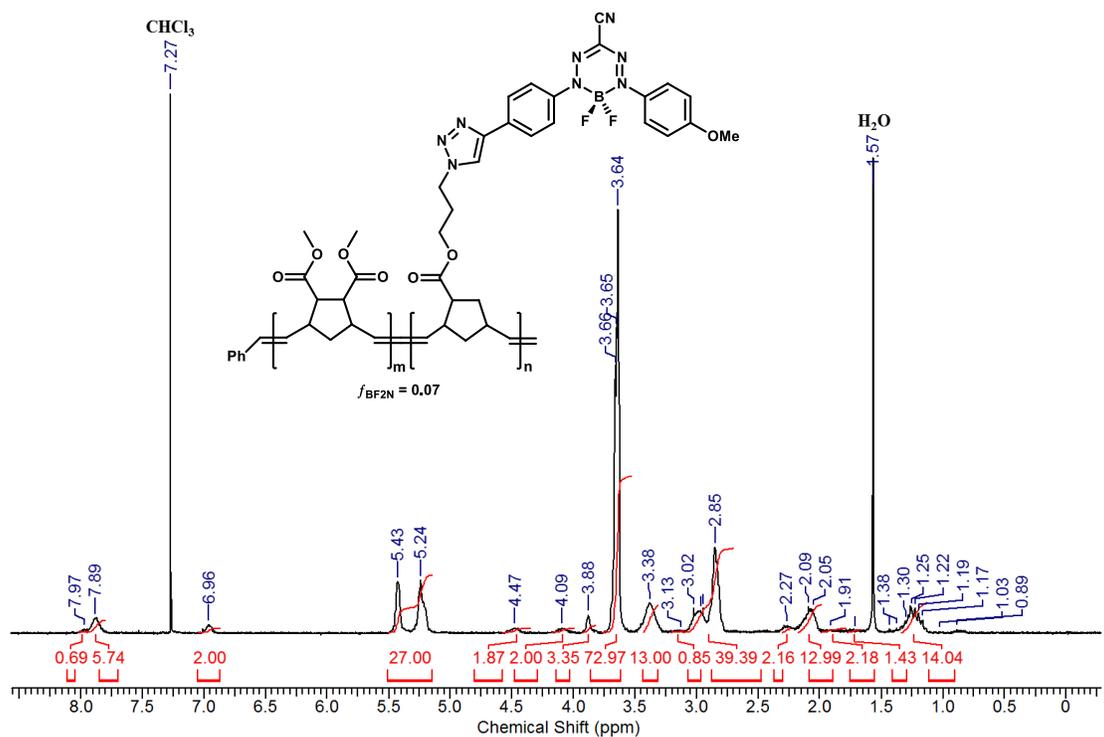


Fig. S14 ^1H NMR spectrum of $(\text{PDND})_m\text{-}b\text{-(PBF}_2\text{N)}_n$ ($f_{\text{BF}_2\text{N}} = 0.07$) in CDCl_3 .

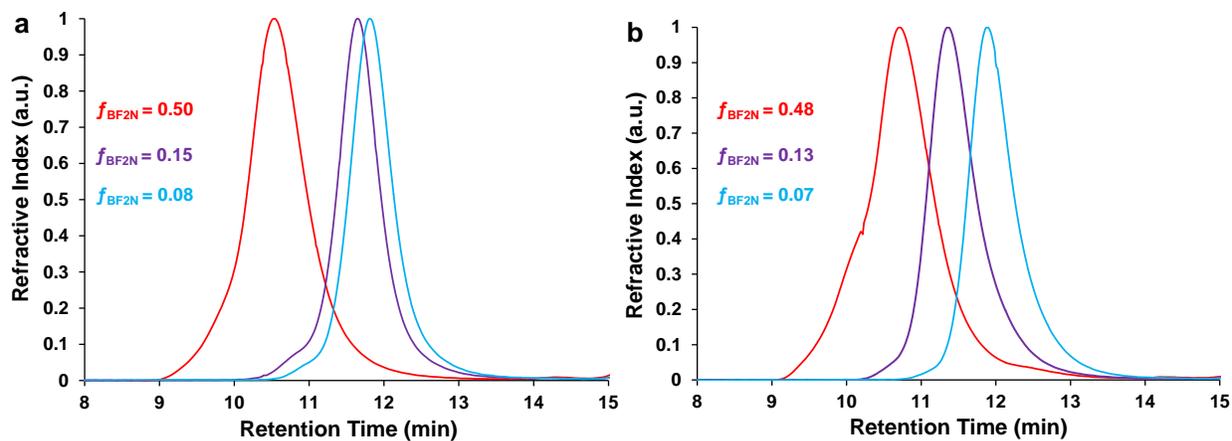


Fig. S15 GPC traces recorded for random copolymers $(\text{PDND})_m\text{-}r\text{-(PBF}_2\text{N)}_n$ (a), and block copolymers $(\text{PDND})_m\text{-}b\text{-(PBF}_2\text{N)}_n$ (b) in DMF.

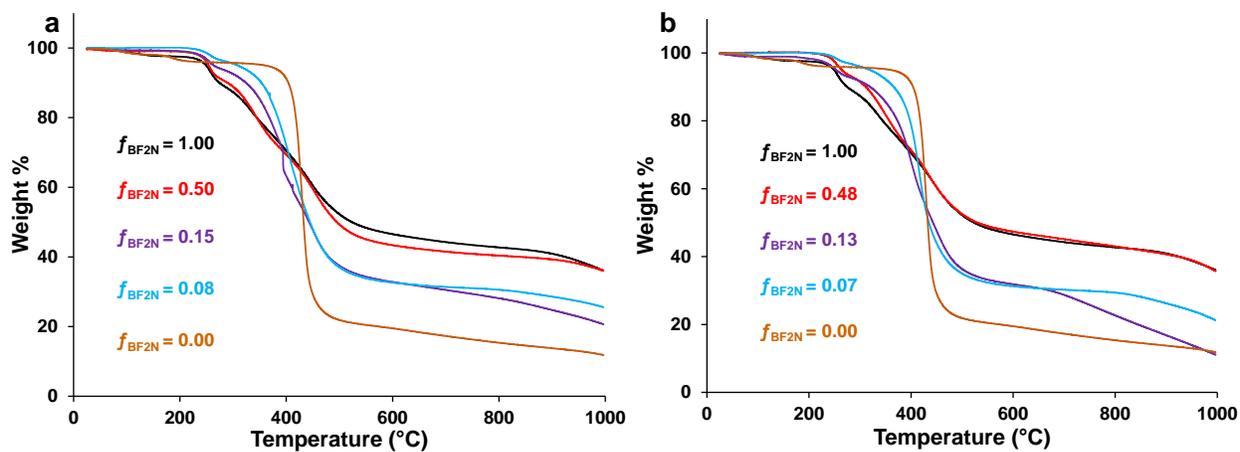


Fig. S16 TGA data recorded for homopolymers **PBF₂N** and **PDND**, random copolymers **(PDND)_m-r-(PBF₂N)_n** (a), and block copolymers **(PDND)_m-b-(PBF₂N)_n** (b) under a N₂ atmosphere.

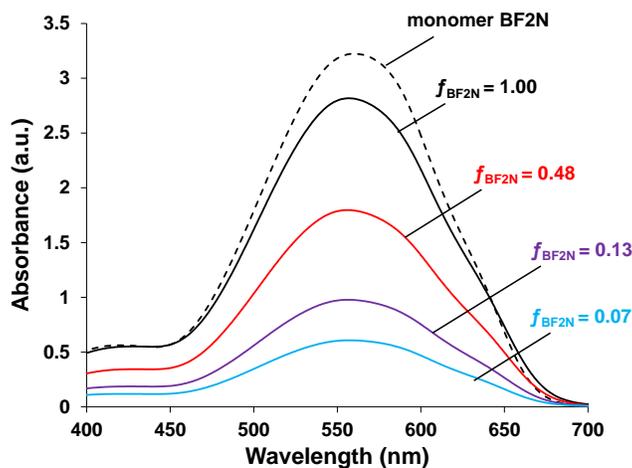


Fig. S17 UV-vis absorption spectra of monomer **BF₂N**, homopolymer **PBF₂N**, and block copolymers **(PDND)_m-b-(PBF₂N)_n** recorded for 0.05 mg mL⁻¹ CH₂Cl₂ solutions.

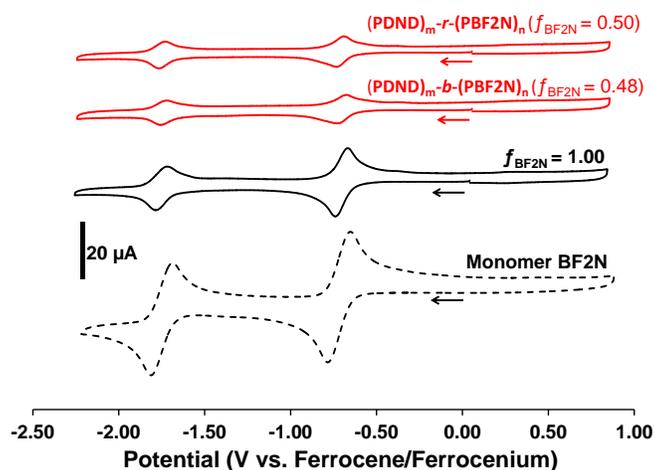


Fig. S18 Cyclic voltammograms of 1 mM (calculated using an average molar mass for blocks and random copolymers) CH_2Cl_2 solutions of monomer BF_2N , homopolymer PBF_2N , random copolymer $(\text{PDND})_m\text{-}r\text{-(PBF}_2\text{N)}_n$ ($f_{\text{BF}_2\text{N}} = 0.50$), and block copolymer $(\text{PDND})_m\text{-}b\text{-(PBF}_2\text{N)}_n$ ($f_{\text{BF}_2\text{N}} = 0.48$) containing 0.1 M $[\text{nBu}_4\text{N}][\text{PF}_6]$ as supporting electrolyte recorded at a scan rate of 250 mV s^{-1} . Voltammograms were referenced internally against the ferrocene/ferrocenium redox couple.