

Supplementary Data

FT-IR

Table S1. Main vibration modes in the PFDPP family of polymers

Vibration mode	PFDPP-1	PFDPP-2	PFDPP-3
	Wave number (cm ⁻¹)		
=C-H st	3066	3066	3066
CH ₃ ν _{as}	2959	2959	2956
CH ₂ ν _{as}	2925	2925	2919
CH ₂ ν _s	2852	2852	2852
C=O st	1664	1658	1658
C-N st	1537	1544	1551
C=C st	1510	1510	1510
δ(C-H) in the thiophene plane	1068	1061	1061
S, thiophene	793	793	786
C-H out of plane of the thiophene ring ν	731	731	732
Deformation C-H out of plane of the thiophene ring	625	632	632

¹H NMR

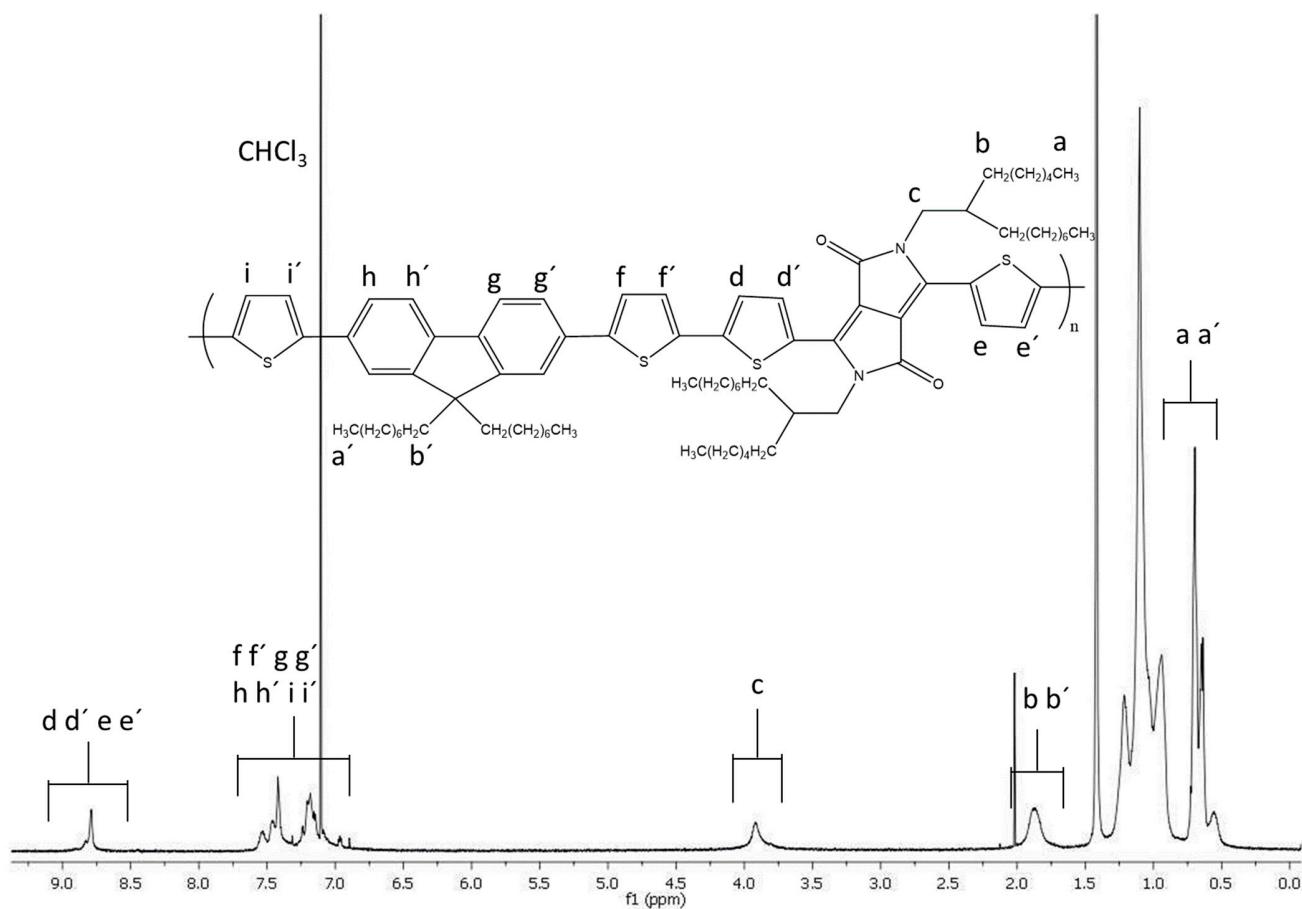


Figure S1. ¹H NMR spectra of PFDPP-2 in CDCl₃

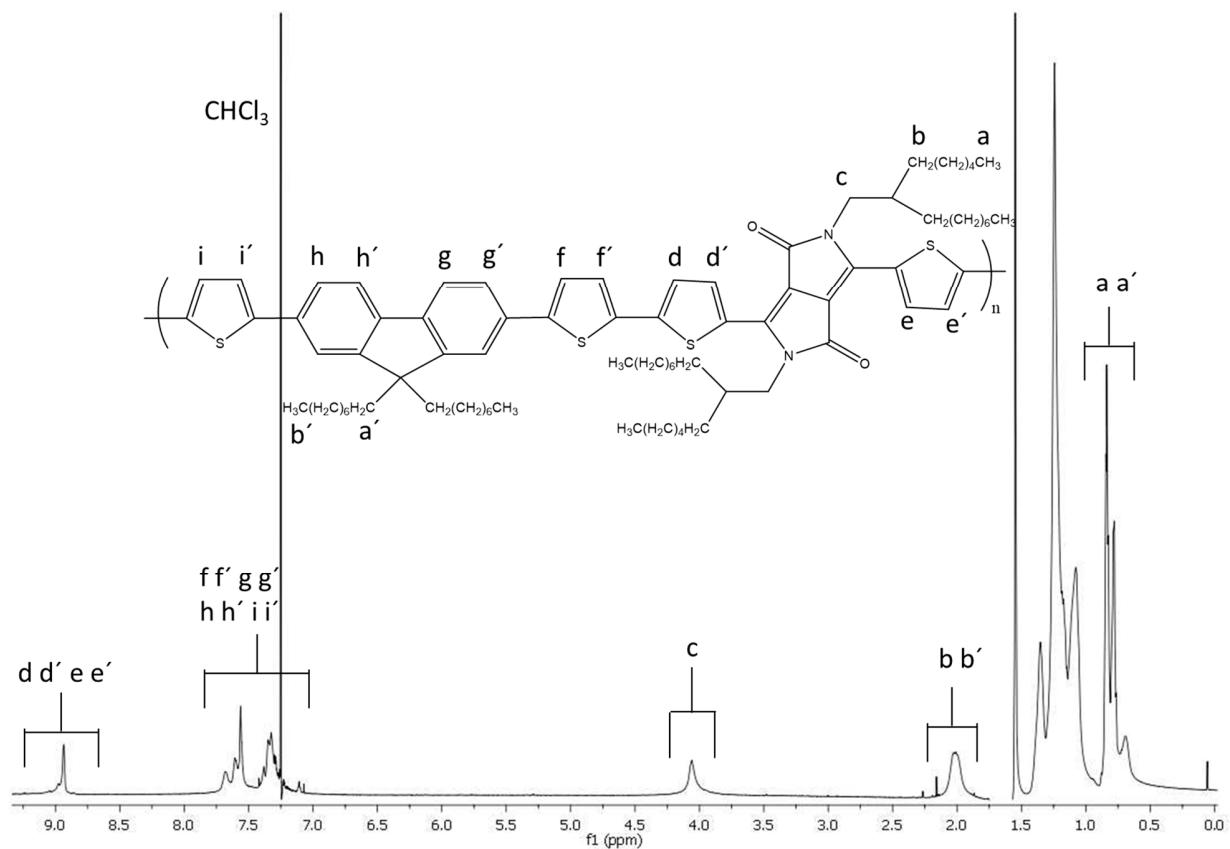


Figure S2. ^1H NMR spectra of PFDPP-3 in CDCl_3

GPC

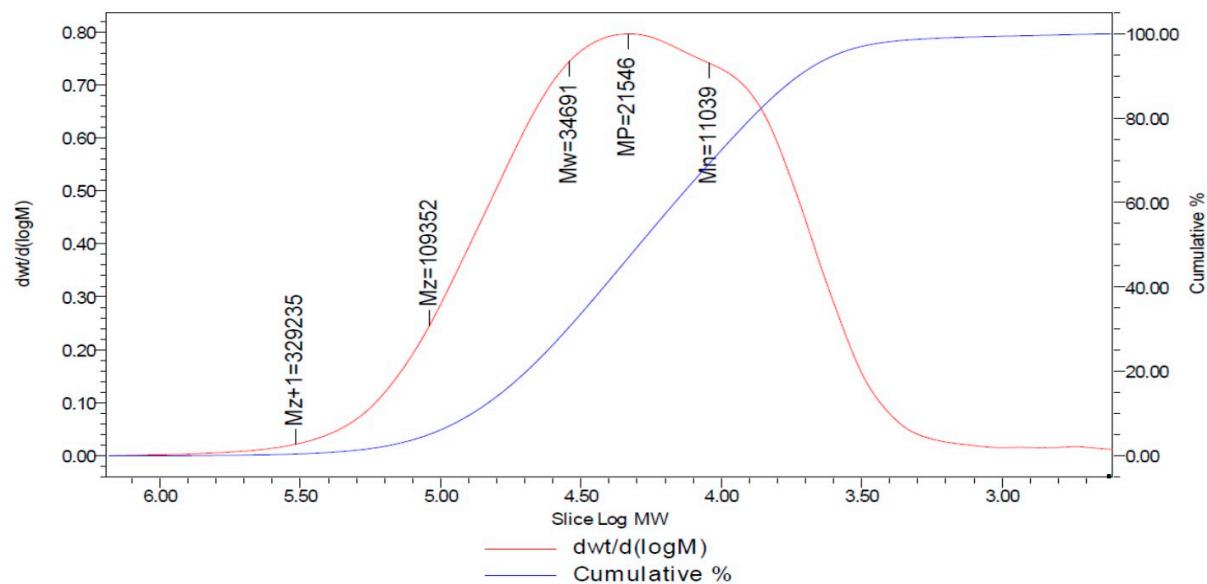


Figure S3. Molecular weight distribution of PFDPP-1

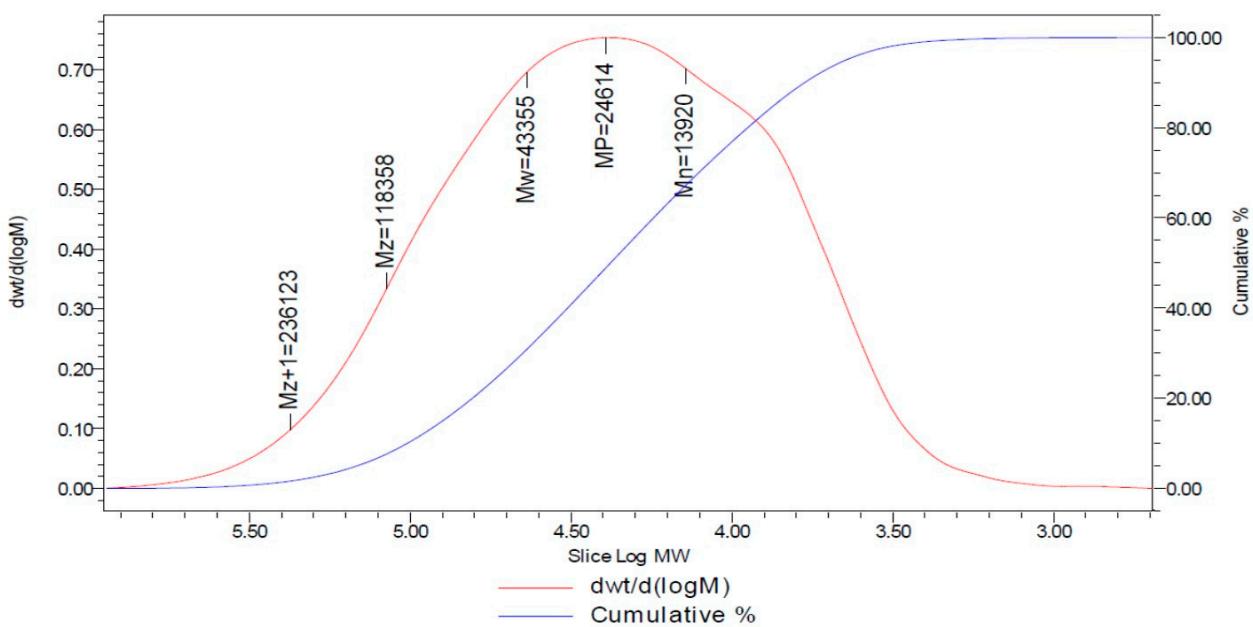


Figure S4. Molecular weight distribution of **PFDPP-2**

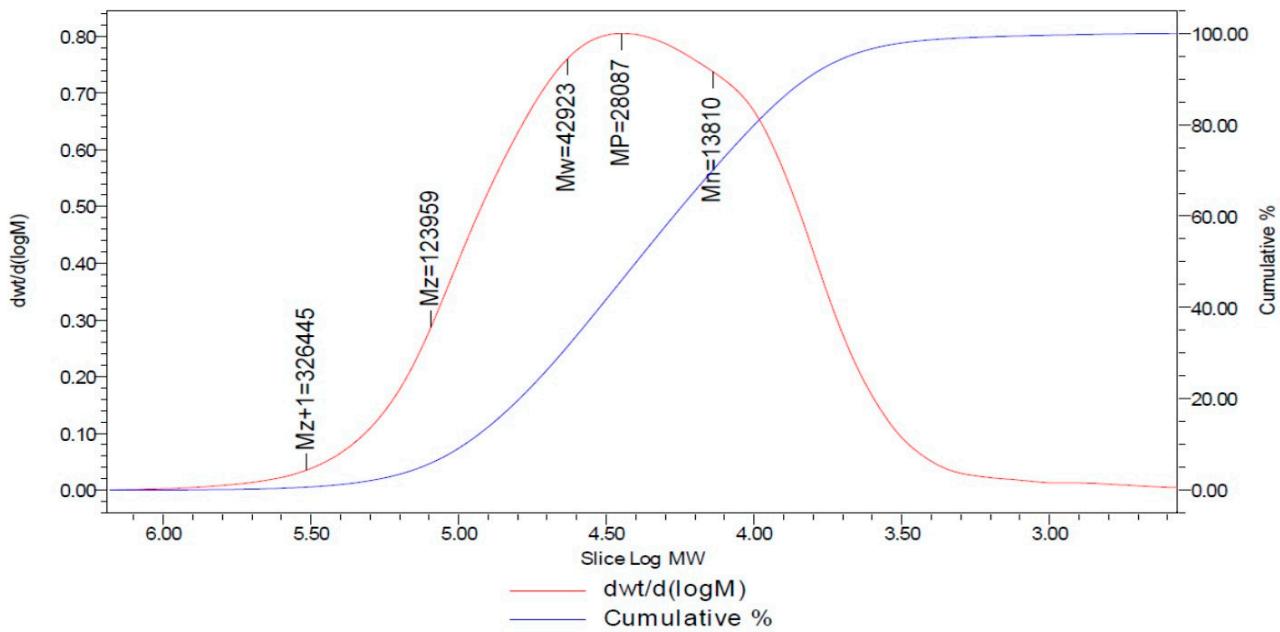


Figure S5. Molecular weight distribution of **PFDPP-3**

Table S2. Fitting of lifetime fluorescence of the PFDPP polymers

Polymer	τ_1 (ns)	B ₁	τ_2 (ns)	B ₂
PFDPP-1	0.78	0.013	-	-
PFDPP-2	0.77	0.029	-	-
PFDPP-3	0.70	0.013	1.16	0.023

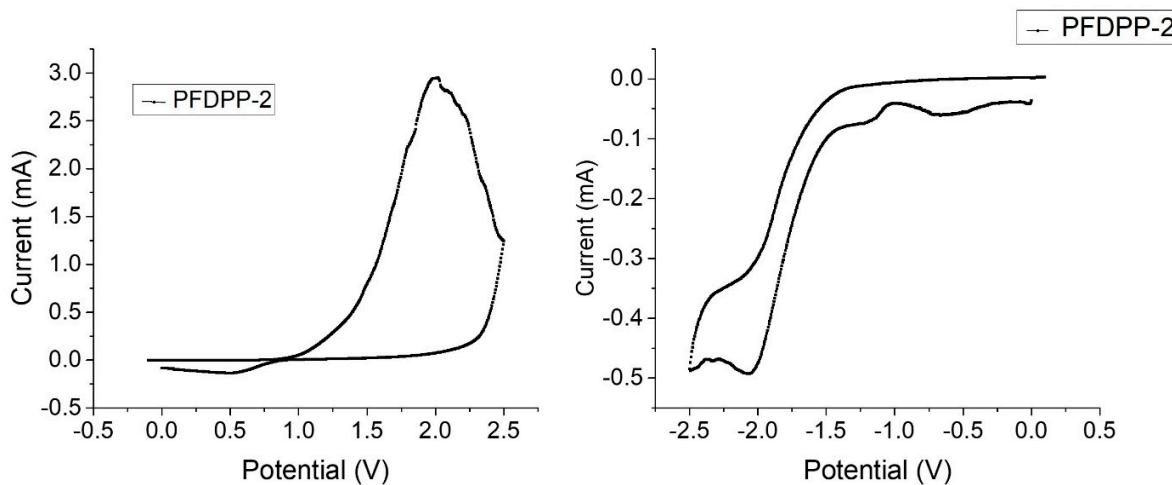


Figure S6. Cyclic voltammograms of PFDPP-2 film polymer (oxidation left and reduction right).

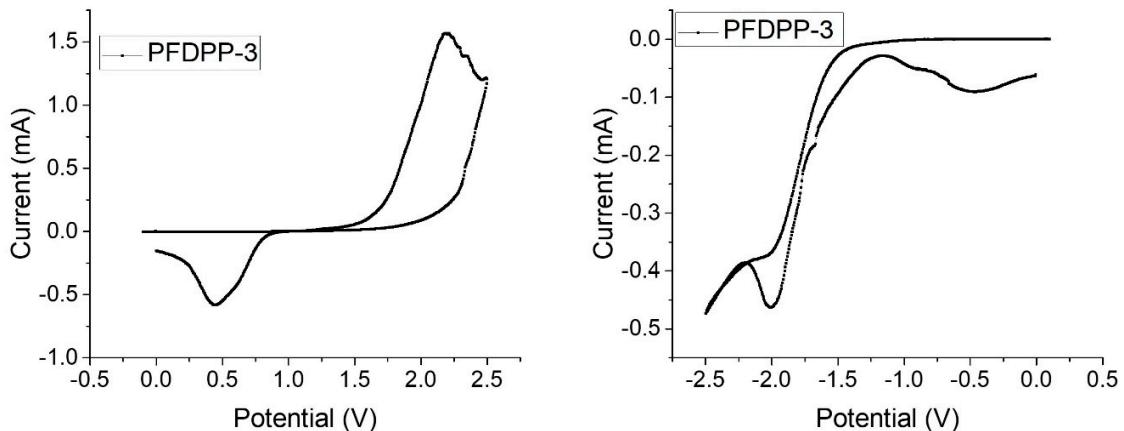


Figure S7. Cyclic voltammograms of PFDPP-3 film polymer (oxidation left and reduction right).