



Polym. Chem.

PAPER

Electronic Supplementary Information for

Received 00th January 20xx,
Accepted 00th January 20xx

DOI: 10.1039/x0xx00000x

www.rsc.org/

Tailor design of renewable copolymers based on poly(1,4-butylene 2,5-furandicarboxylate) and poly(ethylene glycol) with refined thermal properties

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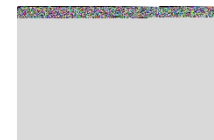
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1 Structural characterisation of the PBF/PEGFs

1.1 FTIR spectra

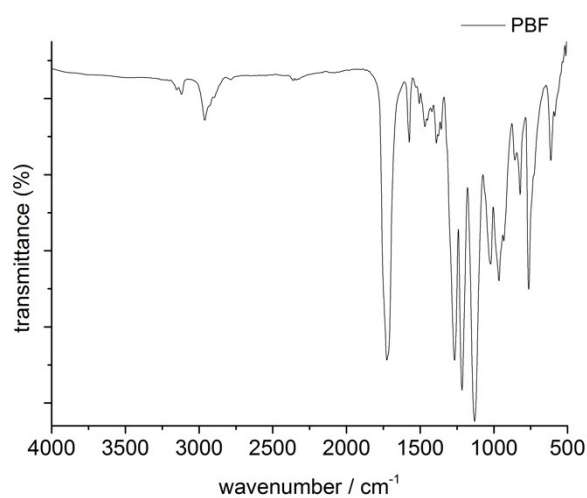


Fig. S1 FTIR spectrum of PBF homopolymer.

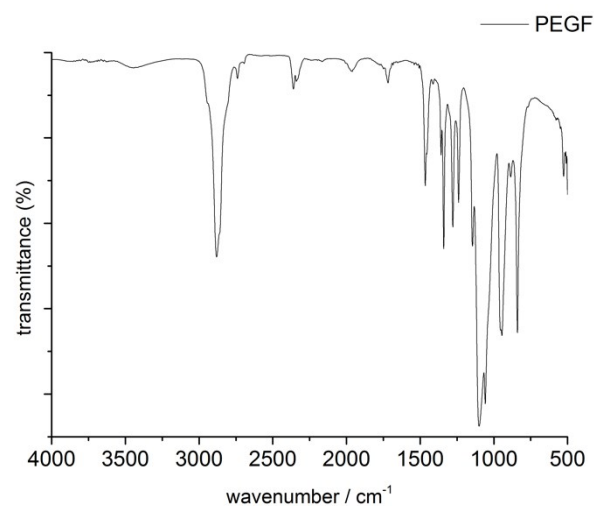


Fig. S2 FTIR spectrum of PEGF homopolymer.

1.2 NMR spectra

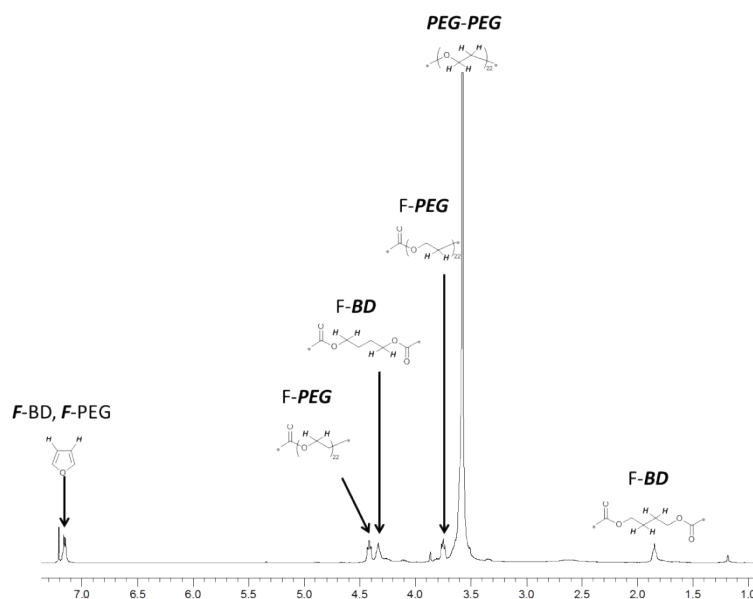


Fig. S3 ^1H NMR spectra of PBF/PEGF2 in deuterated chloroform.

Table S1 Assignments of the main ^{13}C NMR resonances (CDCl_3) of PBF/PEGF copolymers and of PBF and PEGF homopolymers.

assign. ^a	δ / ppm					
	PBF ^b	1 ^b	2	3	4	PEGF ^c
C=O;	158.0	160.5	158.0	157.9	157.9	161.0
F-BD, F-PEG						
C2, C5;	146.8	146.6	146.8	146.7	146.7	154.7
F-BD, F-PEG						
C3, C4;	118.5	119.7	118.5	118.6	118.6	119.0
F-BD, F-PEG						
OCH₂;	-	69.5	70.5	70.6	70.6	70.56
PEG-PEG						
C(O)OCH₂CH₂O;	-	68.6	68.9	68.9	68.9	68.8
F-PEG						
C(O)OCH₂;	64.9	66.7	64.9	64.9	n.d.	-
F-BD						
C(O)OCH₂;	-	66.7	64.5	64.5	64.5	63.0
F-PEG						
C(O)OCH₂CH₂;	25.3	24.4	25.3	25.2	n.d.	-
F-BD						

^a ^{13}C NMR resonances assignment and corresponding group and diad (two adjacent structural units F-BD, F-PEG and/or PEG-PEG). F, BD and PEG stand for 2,5-furandicarboxylate, 1,4-butanediyl and PEG moieties, respectively. ^b ^{13}C NMR spectra in deuterated TFA. ^c ^{13}C NMR spectra were recorded in deuterated chloroform.

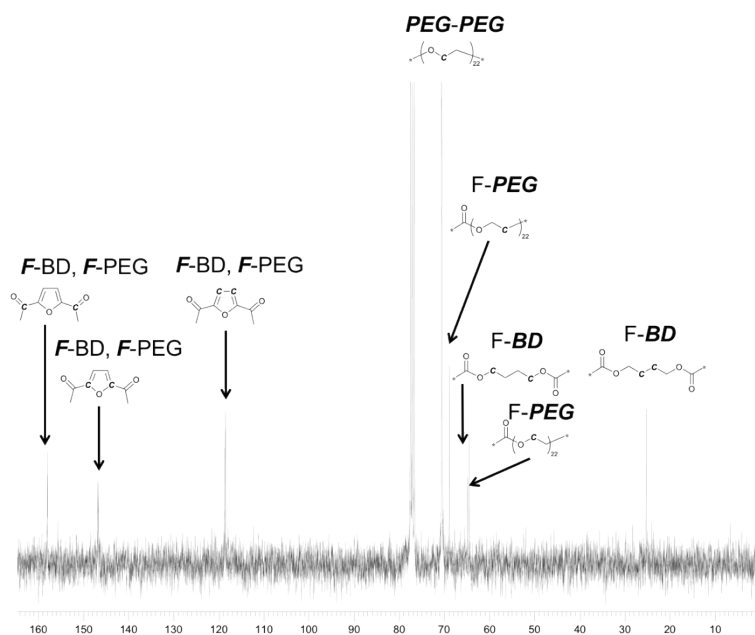


Fig. S4 ^{13}C NMR spectra of PBF/PEGF2 in deuterated chloroform.

2 Thermal behaviour of PBF/PEGFs

2.1 TGA traces

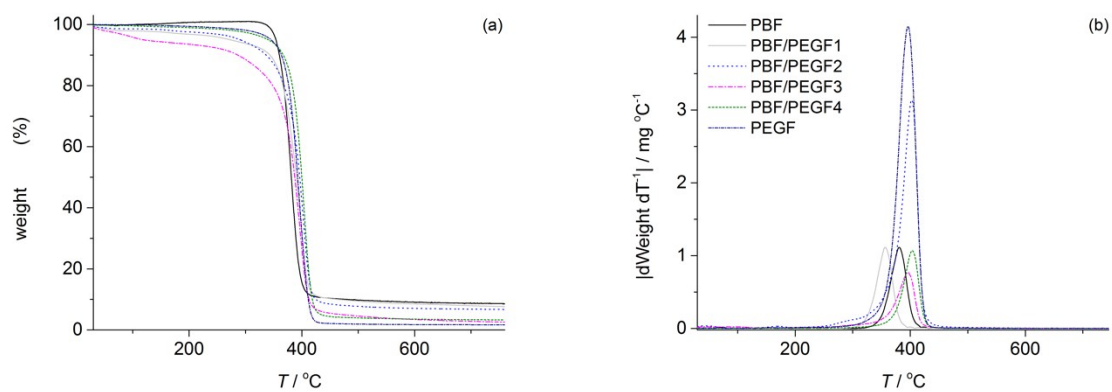


Fig. S5 Thermogravimetric curves of the PBF/PEGFs, PBF and PEGF: (a) TG and (b) DTG.

2.2 DMTA trace

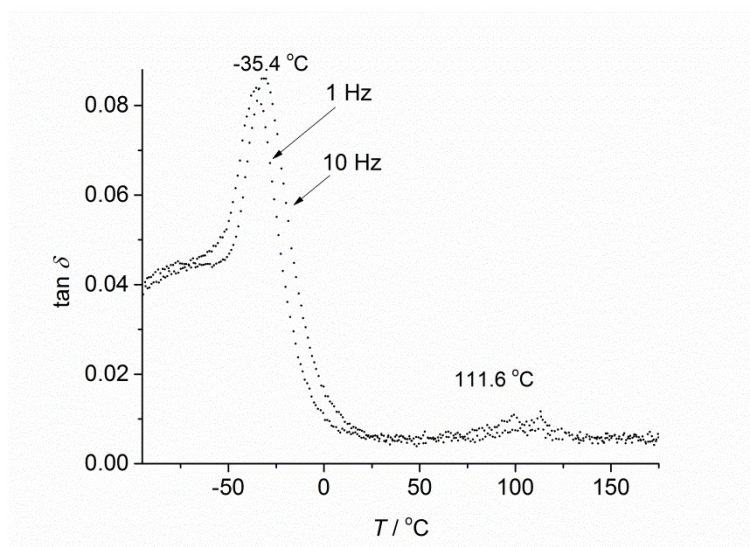


Fig. S6 DMTA traces of PBF/PEGF1 at 1 and 10 Hz.

3 XRD pattern

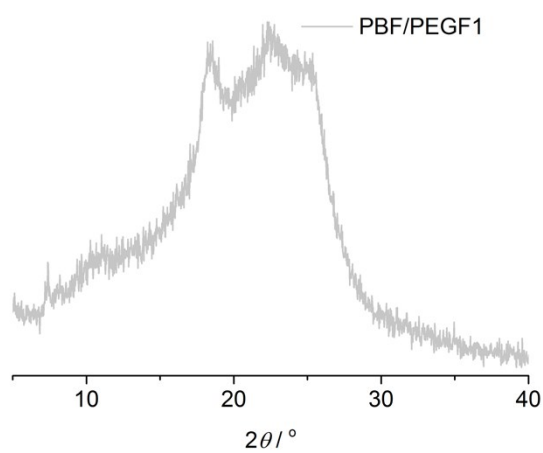


Fig. S7 XRD pattern of PBF/PEGF1.