## **Electronic Supplementary Information**

## Chemical Composition Characterization of Poly(vinylidene fluoridechlorotrifluoroethylene) Based Copolymers with F-H Decoupled <sup>1</sup>H NMR

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## **Result and Discussion**



Fig. S1 <sup>19</sup>F NMR spectrum of the P(VDF-TrFE-CTFE) terpolymer.

Table S1 Chemical shifts and assignments of <sup>19</sup>F NMR peaks of the P(VDF-TrFE-CTFE) terpolymer.

Peak no.	Sequence	Chemistry shift (ppm)
1	-CF <sub>2</sub> CH <sub>2</sub> CF <sub>2</sub> CH <sub>2</sub> CF <sub>2</sub> -	-92.4
2	-CH <sub>2</sub> CH <sub>2</sub> CF <sub>2</sub> CH <sub>2</sub> CF <sub>2</sub> -	-94.5 ~ -96.2
3	-CF2CFClCF2CFClCF2-	-106.7 ~ -109.7
4	$-CF_2CH_2CF_2CF_2CFCI-$	-109.7
5	-CF2CFClCF2CFClCH2-	-112.3 ~ -112.9
6	-CF <sub>2</sub> CH <sub>2</sub> CF <sub>2</sub> CF <sub>2</sub> CFH-	-113.0
7	$-CF_2CH_2CF_2CF_2CH_2-$	-114.7
8	$-CH_2CF_2CF_2CH_2CH_2-$	-117.0
9	-CH <sub>2</sub> CF <sub>2</sub> CF <sub>2</sub> CFClCH <sub>2</sub> -	-119.6 ~ -120.3
10	-CF2CF2CF2CFCICH2CF2-	-121.1 ~ -121.8
11	-CF <sub>2</sub> CFHC <b>F</b> <sub>2</sub> CFHCF <sub>2</sub> -	-122.9
12	-CF <sub>2</sub> CFHC <b>F</b> <sub>2</sub> CFHCH <sub>2</sub> -	-123.6
13	-CH <sub>2</sub> CF <sub>2</sub> CF <sub>2</sub> CFHCH <sub>2</sub> -	-130.5
14	-CF <sub>2</sub> CFClC <b>F</b> <sub>2</sub> CFHCH <sub>2</sub> -	-131.2
15	-CH <sub>2</sub> CF <sub>2</sub> CF <sub>2</sub> CF <sub>2</sub> CH <sub>2</sub> -	-199.0

Fig. S1 shows the <sup>19</sup>F NMR spectrum of a P(VDF-TrFE-CTFE) terpolymer and the assignments of the peak are listed in Table S1.



Fig. S2 <sup>19</sup>F NMR spectrum of the P(VDF-CTFE) copolymer (VDF/CTFE = 91.0/9.0).



Fig. S3 (A) F-decoupled <sup>1</sup>H NMR spectrum of a P(VDF-CTFE) (VDF/CTFE = 91.0/9.0) sample and (B) <sup>1</sup>H NMR of its full hydrogenation product P(VDF-TrFE).

Fig. S2 shows the <sup>19</sup>F NMR spectrum of a P(VDF-CTFE) copolymer. The molar ratio of main chain could be calculated by the equation, VDF/CTFE =  $3(I_1 + I_2 + 3I_3 - I_4)/2[I_5 + I_6 + 2(I_4 - I_3)]$ , the chemical composition of P(VDF-CTFE) is calculated to be VDF/CTFE = 91.0/9.8 mol%. Fig. S3 shows the corresponding F-decoupled <sup>1</sup>H

NMR spectra of P(VDF-CTFE) and <sup>1</sup>H NMR of its full hydrogenation product. From Fig. S3B, the peaks integral of  $I_1$ ,  $I_2$ ,  $I_3$  and the equation, TrFE/VDF =  $2I_3/(I_1 + I_2)$ , the chemical composition of P(VDF-CTFE) is calculated to be VDF/CTFE = 91.0/9.0 mol%. From Fig. S3A, the peaks integral of  $I_1$ ,  $I_2$ ,  $I_4$  and the equation, CTFE/VDF =  $I_4/(I_1 + I_2)$ , the chemical composition of the same sample is calculated to be VDF/CTFE = 91.0/8.3 mol%, which is slightly smaller than the other two methods mentioned above. And the reason has been explained in the main text. The composition results calculated from both hydrogenation method and F-H decoupled method agree very well suggesting the accuracy of F-H decoupled <sup>1</sup>H NMR.