Supplementary Information for "Quantitative rationalization of unexpectedly moderate water wettability on poly (vinyl alcohol) surfaces: Thermodynamic Evaluation and Prediction of Surface Hydrogen Bonding"

Zhuohuan Guo, Zhuoyuan Ma * and Dayang Wang *

State Key Laboratory for Inorganic Synthesis and Preparative Chemistry, College of Chemistry, Jilin University, 130012, Changchun, China

* Corresponding authors. E-mail: wangdayang@jlu.edu.cn, mzyjlu@163.com

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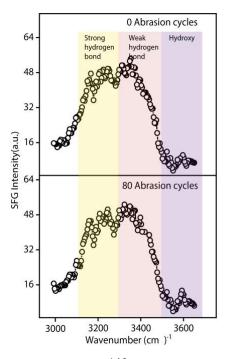
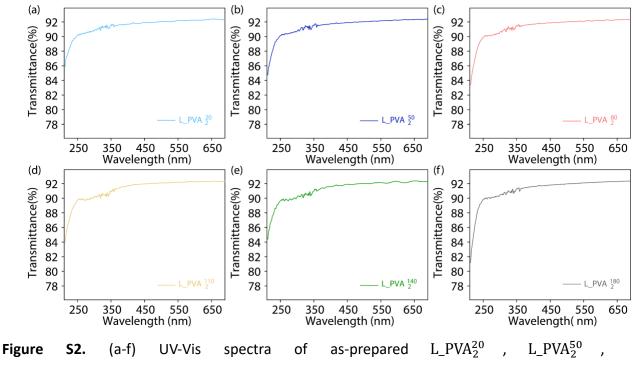


Figure S1. SFG spectra of as-prepared $L_PVA_2^{140}$ thin films before and after 80 cycles of abrasion.



 $L_PVA_2^{80}$, $L_PVA_2^{110}$, $L_PVA_2^{140}$ and $L_PVA_2^{180}$ thin films in the wavelength range of 200-700 nm, respectively.

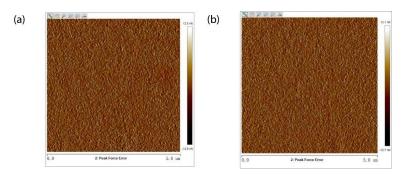


Figure S3. AFM images of as-prepared $L_PVA_2^{20}$ (a) and $L_PVA_2^{180}$ (b) thin films.

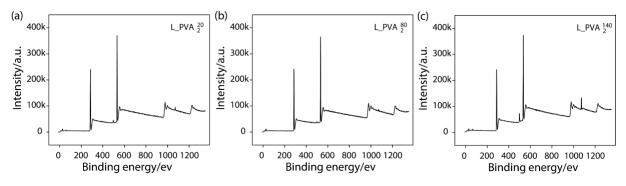


Figure S4. High-resolution XPS spectra of the survey signals of $L_PVA_2^{20}$ (a), $L_PVA_2^{80}$ (b), and

 $L_PVA_2^{140}$ thin films (c).

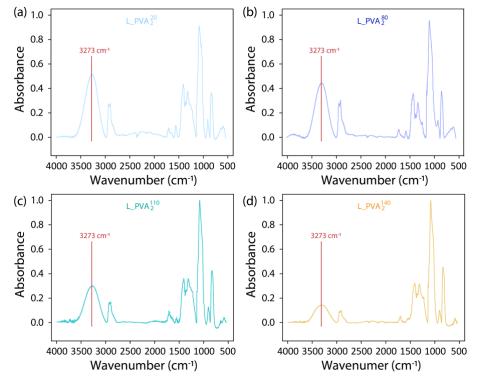


Figure S5. (a-d) ATR-FTIR spectra of as-prepared $L_PVA_2^{20}$, $L_PVA_2^{80}$, $L_PVA_2^{110}$, and $L_PVA_2^{140}$ thin films in the range of 4000-400 cm⁻¹, respectively.

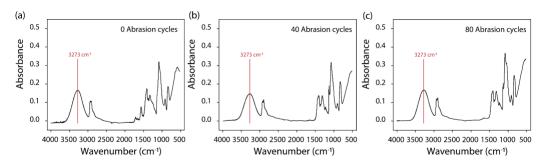


Figure S6. (a-c) ATR-FTIR spectra of as-prepared $L_PVA_2^{140}$ thin films after their surfaces are subjected to 0, 40 and 80 cycles of abrasion, respectively.



Figure S7. (a-c) SEM images of as-prepared $H_PVA_2^{80}$, $L_PVA_7^{RT}$, and $M_PVA_G^{RT}$ films, respectively.

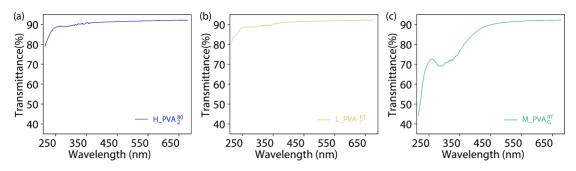


Figure S8. (a-c) UV-Vis spectra of as-prepared $H_PVA_2^{80}$, $L_PVA_7^{RT}$, and $M_PVA_G^{RT}$ films in the wavelength range of 200-700 nm, respectively.

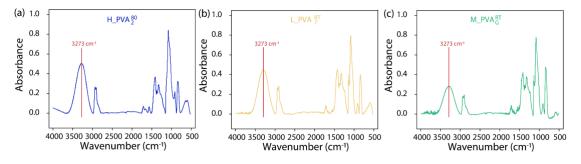


Figure S9. (a-c) ATR-FTIR spectra of $H_PVA_2^{80}$, $L_PVA_7^{RT}$, and $M_PVA_G^{RT}$ films in the range of 4000-400 cm⁻¹, respectively.

Table S1. Summary of the fitting parameters of ATR-FTIR spectra on $L_PVA_2^T$, $H_PVA_2^{80}$, $L_PVA_7^{RT}$ and $M_PVA_G^{RT}$ thin films.

$L_PVA_2^T$	OH_t group				OH_g group				
	peak position	Standard errors	peak area	Standard errors	peak positi on	Standard errors	peak area	Standard errors	R ²
L_PVA ₂ ²⁰	3222 cm ⁻¹	2.4 cm ⁻¹	59.5	3.0	3356	3.3 cm ⁻¹	41.7	3.0	0.9937
L_PVA ₂ ⁸⁰	3222 cm ⁻¹	1.9 cm ⁻¹	50.5	1.9	cm ⁻¹ 3356 cm ⁻¹	2.2 cm ⁻¹	39.6	1.9	0.9981
$L_PVA_2^{110}$	3222 cm ⁻¹	2.3 cm ⁻¹	37.1	1.5	3356	2.7 cm ⁻¹	30.0	1.5	0.9951
L_PVA ₂ ¹⁴⁰	3222 cm ⁻¹	2.0 cm ⁻¹	22.7	0.6	cm ⁻¹ 3356 cm ⁻¹	2.2 cm ⁻¹	42.5	0.6	0.9959
H_PVA ₂ ⁸⁰	3222 cm ⁻¹	2.6 cm ⁻¹	56.5	3.8	3356	3.0 cm ⁻¹	54.0	4.6	0.9962
L_PVA ₇ ^{RT}	3222 cm ⁻¹	5.0 cm ⁻¹	49.8	4.7	cm ⁻¹ 3356 cm ⁻¹	4.0 cm ⁻¹	46.6	4.8	0.9953
M_PVA _G ^{RT}	3222 cm ⁻¹	2.7 cm ⁻¹	28.5	3.7	3356 cm ⁻¹	3.5 cm ⁻¹	24.7	2.7	0.9970

Bootstrapping Validation of $\gamma_t^{s,p*}$ and $\gamma_g^{s,p*}$ values.

Based on the determination method for $\gamma_t^{s,p*}$ and $\gamma_g^{s,p*}$ values reported in Section 3.3, Line 24, and the data listed in Table 1 (at least three sets), we applied the multivariable linear regression model in Origin to fit Equation 13 and obtained the following values that satisfy the condition: 0 mN/m < $\gamma_t^{s,p*}$ and $\gamma_g^{s,p*} < 50$ mN/m:

For
$$\gamma_t^{s,p*}$$
: $(x_1 = 6.2 \ mN/m, x_2 = 8.4 \ mN/m, x_3 = 8.0 \ mN/m)$
For $\gamma_g^{s,p*}$: $(x_1 = 9.0 \ mN/m, x_2 = 8.5 \ mN/m, x_3 = 9.8 \ mN/m)$

The 95% confidence intervals for $\gamma_t^{s,p*}$ and $\gamma_g^{s,p*}$ were calculated using Equation (13), resulting in the following: $\gamma_t^{s,p*}$ 95% CI: 7.5 ± 1.5 mN/m; $\gamma_g^{s,p*}$ 95% CI: 9.1 ± 1.2 mN/m. The confidence interval formula is described as follows:

$$CI = \bar{x} \pm t \times \frac{SD}{\sqrt{n}} \tag{S1}$$

where \bar{x} is the sample mean, t is the critical value from the t-distribution (2.3534), SD is the standard deviation, and n is the sample size.