

Supplementary materials

**Effective removal of Pb(II) from synthetic wastewater
using $\text{Ti}_3\text{C}_2\text{T}_x$ MXene**

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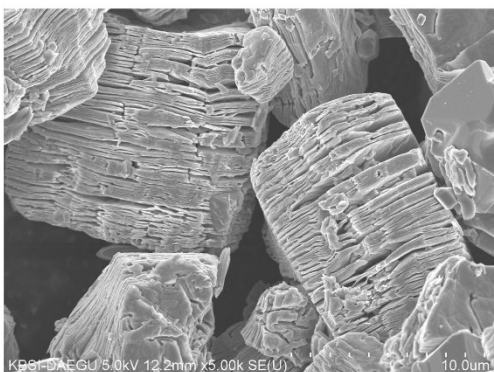
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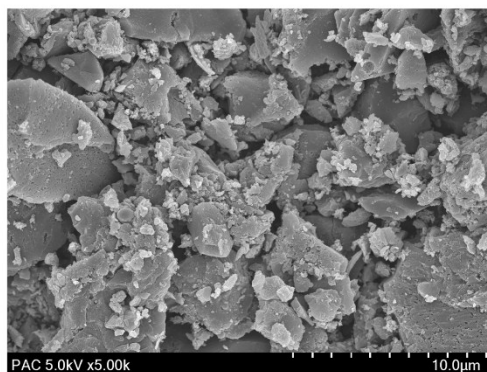
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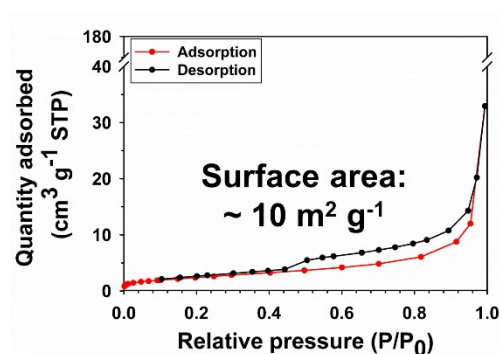
(a.1) SEM



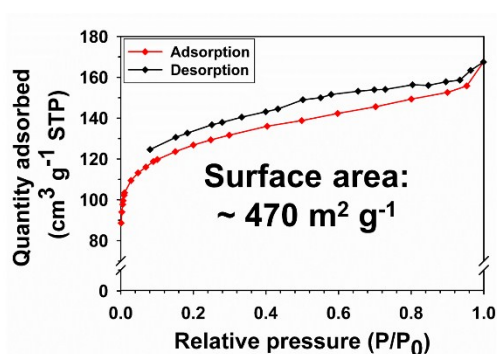
(b.1) SEM



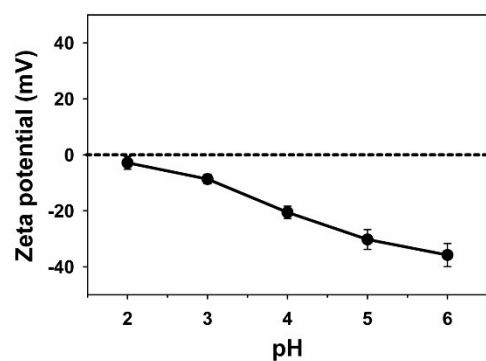
(a.2) BET



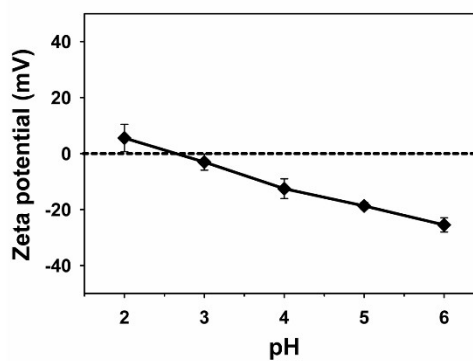
(b.2) BET



(a.3) Zeta potential

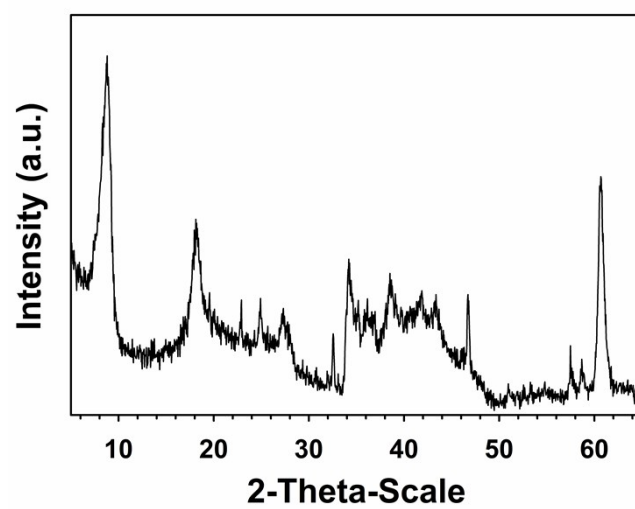


(b.3) Zeta potential



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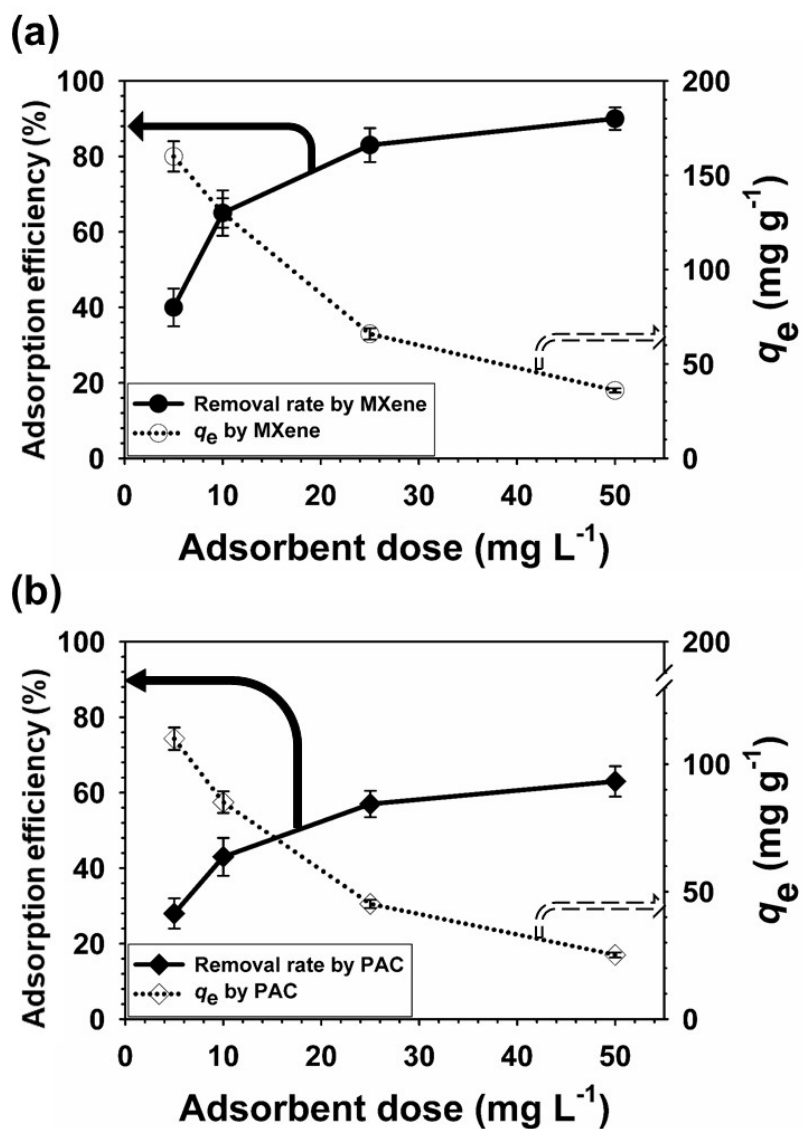
21 **Fig. S1.** Various characterizations of (a) MXene and (b) PAC using (1) SEM, (2) BET, and (3)
22 zeta potential.



24

25 **Fig. S2.** XRD analysis of MXene to confirm crystallinity.

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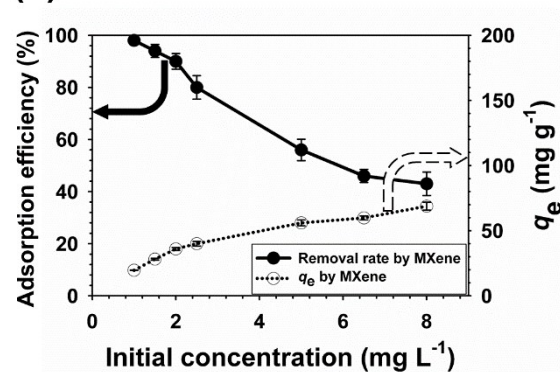
28

29 **Fig. S3.** Influence of (a) MXene and (b) PAC dosage on the q_e and adsorption efficiency of
 30 Pb(II) (Conditions: Pb^{2+} concentration = 2 mg L^{-1} ; temperature = 293 K ; pH = 6; contact time
 31 = 2 h).

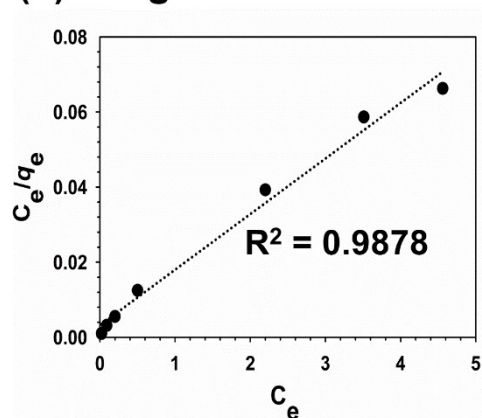
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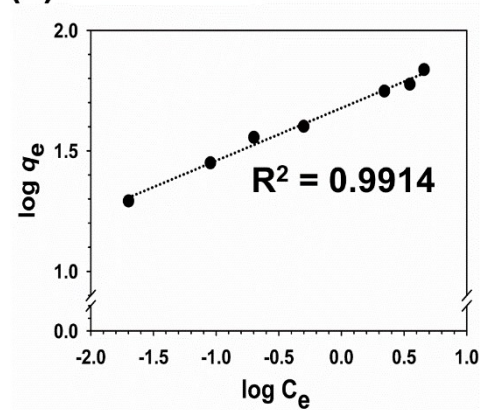
(a) Isotherm



(b) Langmuir



(c) Freundlich

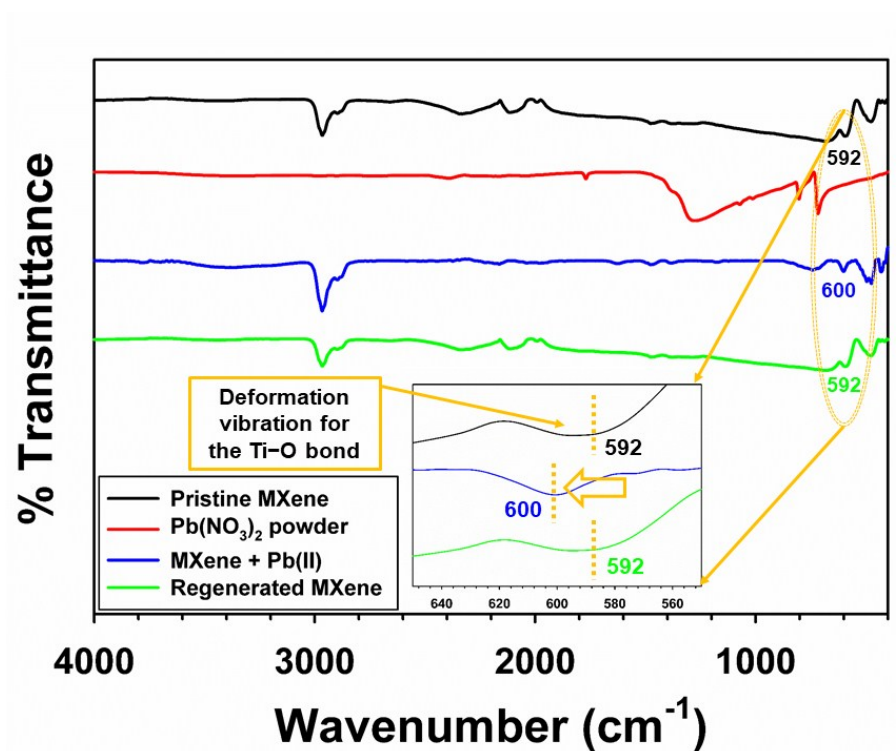


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35 **Fig. S4.** (a) Influence of initial concentration on the q_e and adsorption efficiency of Pb(II). (b)
 36 Langmuir and (c) Freundlich isotherm model to clarify isotherm study of Pb(II) adsorption by
 37 MXene (Conditions: MXene dose = 50 mg L⁻¹; Pb²⁺ concentration = 1 to 8 mg L⁻¹; temperature
 38 = 293 K; pH = 6; contact time = 2 h).

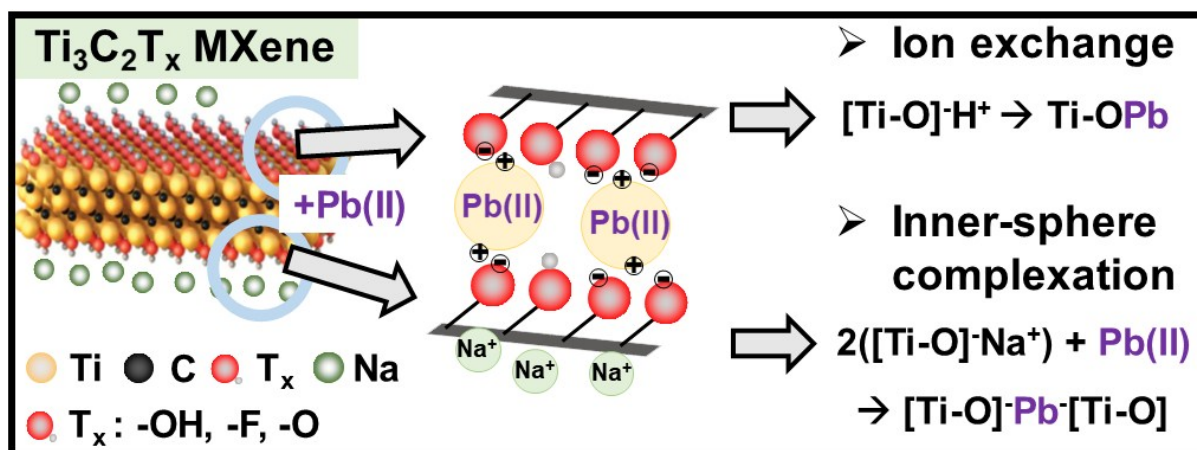
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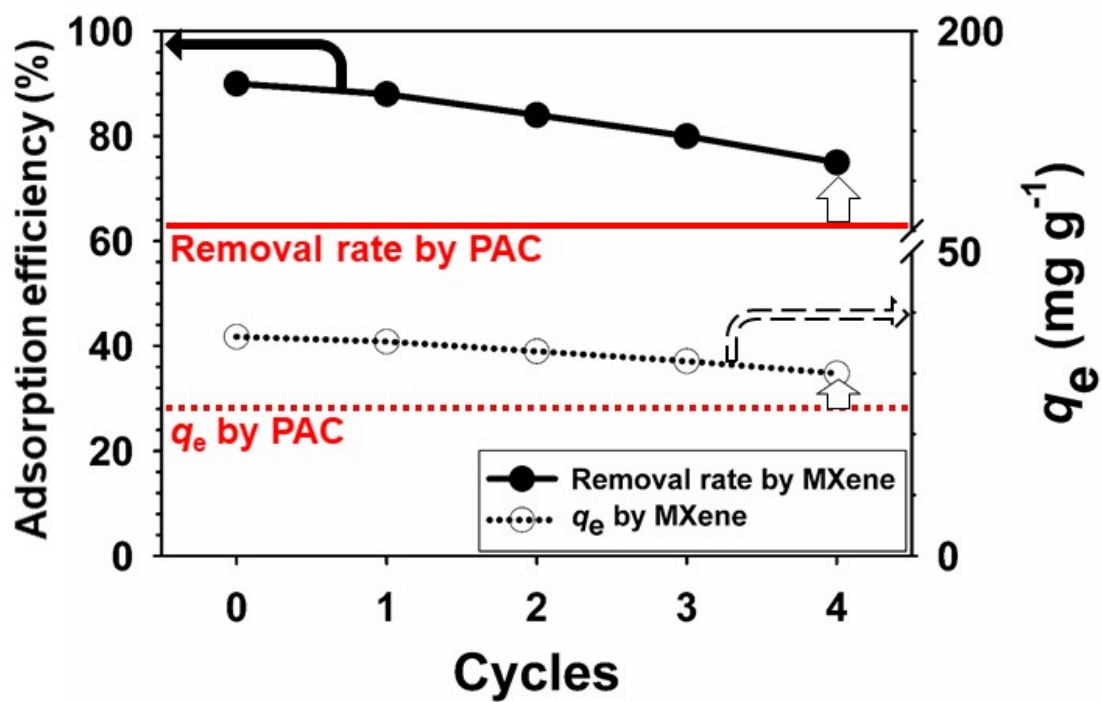
41

42 **Fig. S5.** Confirmation of peak (Ti-O) shift after adsorption of Pb(II) on MXene using FTIR
 43 technique.



45

46 **Fig. S6.** Schematic diagram related to Pb(II) adsorption mechanism.



48

49 **Fig. S7.** Confirmation of re-usability for Pb(II) adsorption in four adsorption-desorption cycles
 50 (Conditions: MXene dose = 50 mg L⁻¹; Pb²⁺ concentration = 2 mg L⁻¹; temperature = 293 K;
 51 pH = 6; contact time = 2 h).

52

54 **Table S1.** Kinetic parameters of Pb(II) adsorption by MXene through four different kinetic
 55 models.

Elovich	
a	3.65
b	18.8
R^2	0.521
Intra-particle diffusion	
k_i (mg g min ^{-0.5})	0.750
C	26.9
R^2	0.340
Pseudo-first-order	
q_e (mg g ⁻¹)	3.59
k_1 (min ⁻¹)	0.0097
R^2	0.473
Pseudo-second-order	
q_e (mg g ⁻¹)	36.6
k_2 (g mg-min ⁻¹)	0.0088
R^2	0.999

57 **Table S2.** Thermodynamic parameters for Pb(II) adsorption by MXene.

Thermodynamic parameters				
ΔG (kJ mol ⁻¹) at different temperature (K)			ΔH (kJ mol ⁻¹)	ΔS (kJ mol ⁻¹ K ⁻¹)
293	303	313		
-12.65	-13.88	-14.73	17.82	0.10

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