

Supplementary Materials

Table S1. Synthesis details of 2D layered MXenes from MAX phases.

S. No	Synthesis Process (Ball Milling+ Chemical Etching)	Abbreviation
1	6 h Ball Milled MAX phase	BMAX-6H
2	6 h Ball milled and 12 h etched and dried at 60°C	BM-12H
3	12h etched and dried at 60°C	M-12H
4	24h etched and dried at 60°C	M-24H
5	48h etched and dried at 60°C	M-48H
<i>Etching By HF.</i>		

Calculation of electrochemical parameters:

The electrochemical parameters, such as the specific capacitance (C_s , Fg⁻¹), Coulombic efficiency (η), energy (E), and power density (P) are important parameters for the investigation of the capacitive behavior of electrochemical cells.

Calculation of specific capacitance for three electrode system:

In order to evaluate the performances of the cells, the specific capacitances for the electrodes were tested in three-electrode system. The specific gravimetric capacitance (Fg⁻¹) for the three-electrode setup was calculated from galvanostatic charging and discharging (GCD) curves using the equation 1:

$$C_{sc}(F\ g^{-1}) = \frac{I\Delta t}{\Delta V m} \quad (1)$$

where, I is the discharge (or charge) current, Δt is the discharge (or charge) time, ΔV is the potential window of cycling, and m is the mass loaded on carbon cloth substrate.

Specific capacitance for two electrode cell:

The gravimetric capacitance for symmetric supercapacitor was calculated from galvanostatic charge/discharge curves using the equation 2:

$$CSC(F\ g^{-1}) = \frac{2I\Delta t}{\Delta V m} \quad (2)$$

where Csc (F g⁻¹) is the gravimetric specific capacitance, i (A) is the discharge current, Δt (s) is the discharge time, ΔV is the potential window (excluding the IR drop), and m in gram is the mass of the active material.

Power and energy density calculations:

The power density (P) of the symmetric supercapacitor was calculated from the energy density (E) and the discharge time (t) according to the Equations (3) and (4)

$$E_{cell} = \frac{1}{2} C_{sc} \Delta V^2 \cdot \frac{1}{4 * 3.6} \quad (3)$$

$$P_{cell} = \frac{E_{cell}}{\Delta t} \quad (4)$$

where E is the energy density (Wh kg^{-1}), C is the measured device capacitance (Fg^{-1}) calculated using eq (2), V is the potential window (V), P is the power density (W kg^{-1}), and Δt is the discharge (or charge) time (s) from the CD curves.

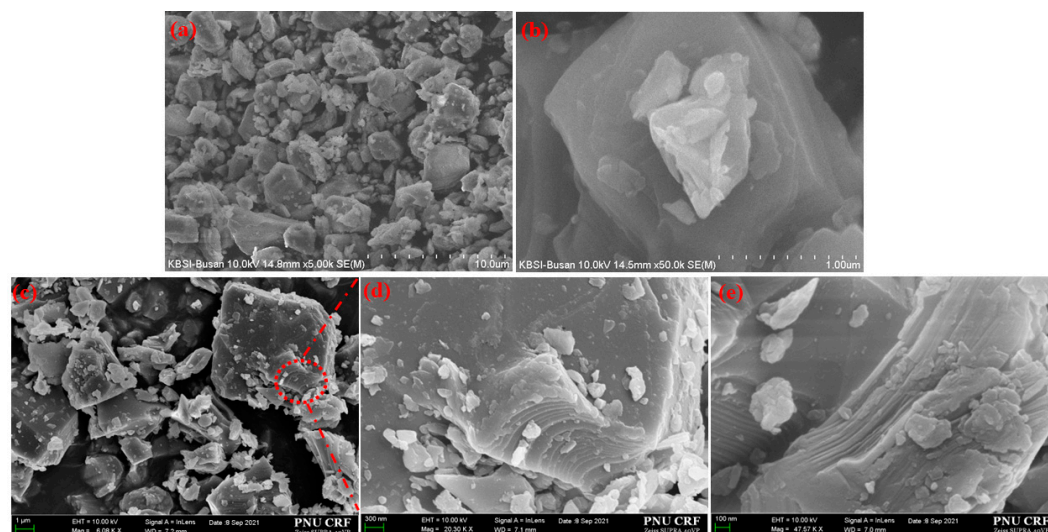


Figure S1. FESEM images of (a and b) pristine Ti_3AlC_2 , (c, d, e) Ball milled MAX phase at low and high magnification.

Table S2. The comparison of MXene electrode supercapacitor's various parameters in 3 electrode system with aqueous electrolyte.

No.	Material	Electrolyte	Potential Window(V)	Specific Capacitance (F g^{-1})	Ref
1	Ti_3C_2	1M NaCl	1.2	127.6 at 0.1 A g^{-1}	[1]
2	$\text{Ti}_3\text{C}_2/\text{TiO}_2$ -nanowire	6M KOH	0.6	135 at 5 mV/s	[2]
3	Ti_3C_2	0.5M NaCl	1.2	132 at 0.1 A g^{-1}	[3]
4	Ti_3C_2	1M NaCl	0.9	75.76 at 0.1 A g^{-1}	[4]
5	Ti_3C_2	1M H_2SO_4	0.5	105.8 at 1 A g^{-1}	[5]
6	Ti_3C_2	3M KOH	0.8	141.45 at 0.4 A g^{-1}	[6]
7	Ti_3C_2	3M KOH	0.8	146.25 at 0.5 A g^{-1}	This work

Table S3. Surface Area, Pore Size, and Pore Volume Data.

MXene Samples	SSA ($\text{m}^2 \text{ g}^{-1}$) ^a	Pore Volume ($\text{cm}^3 \text{ g}^{-1}$) ^a	Pore diameter (nm) ^a
BMAX-6H	8.57	0.02	-
BM-12H	32.8	0.24	7.6
M-24H	23.7	0.13	22.1
M-48H	17.28	0.08	24.8

Table S4. The impedance parameters derived by equivalent circuit model.

Electrode	$R_s(\Omega)$	$R_{SEI}(\Omega)$	$Q_1(\mu F)$	$R_{ct}(\Omega)$	$Q_2(\mu F)$	$Z_w(\Omega s^{-1/2})$
M-24H	0.64	0.56	23.23	0.43	335	1.63
M-48H	0.74	8.68	5670	0.65	24.23	9.12
BM-12H	0.7	0.40	27.01	0.33	668	1.31

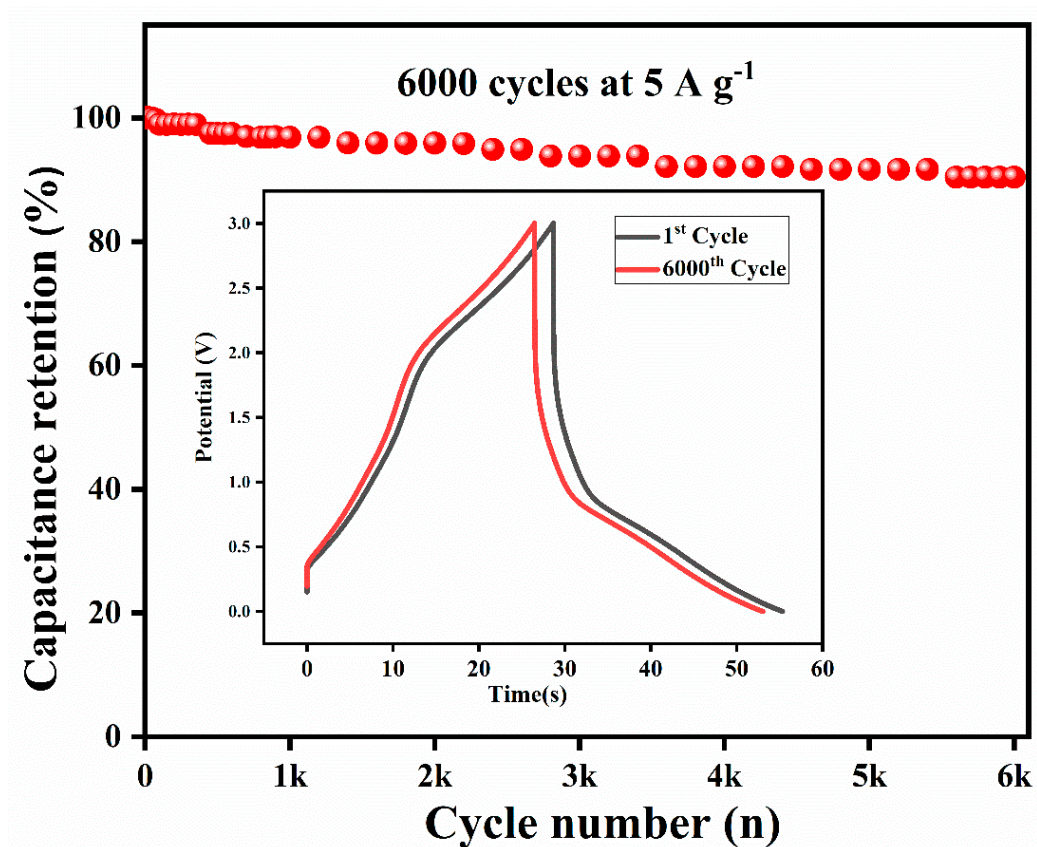


Figure S2. Cyclic stability of MXene coin cell at current density of 5 A g^{-1} .

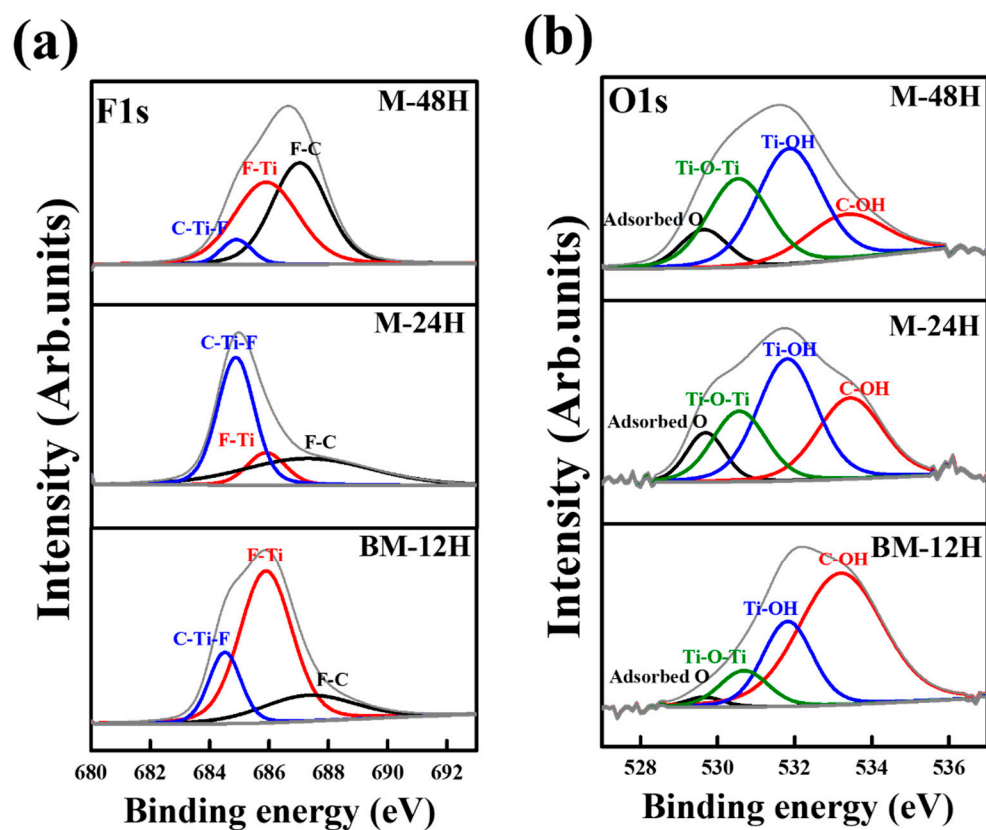


Figure S3. XPS fitting of (a)F1s and (b)O1s for M-48H, M-24, BM-12H.

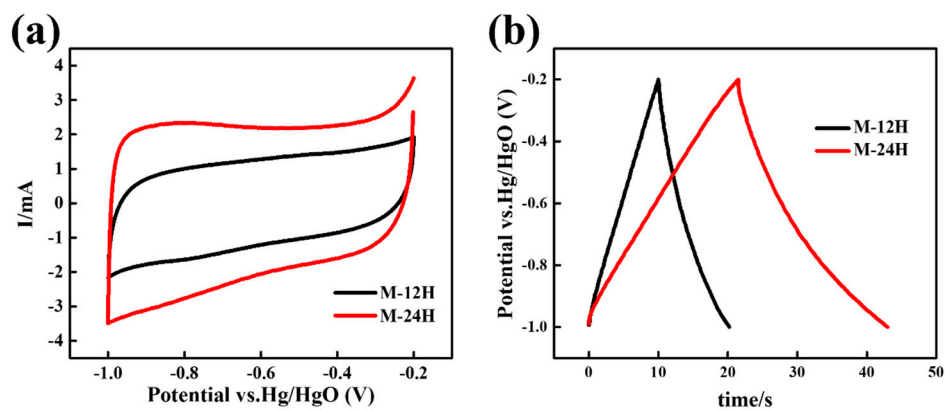


Figure S4. Comparison in (a) CV and (b) GCD for M-12H and M-24H.

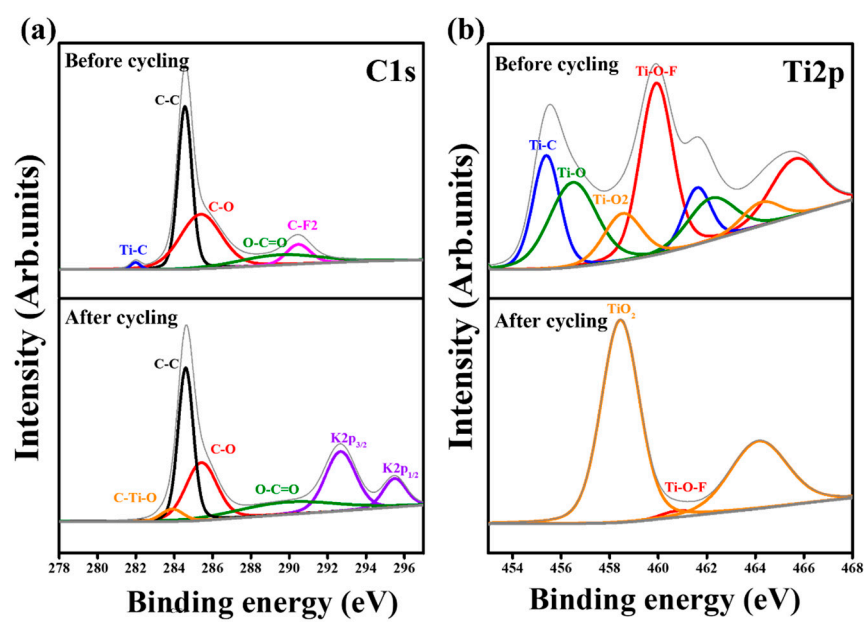


Figure S5. Comparison of XPS spectra in (a)C1s and (b)Ti2p before and after cycling tested BM-12H electrode.

References

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