



## Supplementary Material :Temperature-Dependent Electrical Properties of Al<sub>2</sub>O<sub>3</sub> Passivated Multilayer MoS<sub>2</sub> Thin-Film Transistors

**Keywords:** transition metal dichalcogenide; molybdenum disulfide; thin-film transistor; passivation; contact resistance; intrinsic mobility



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Atomic %	0-К	AI-K	S-KA	Mo-KA	Si-KA
1	57.83	41.18	0.26	0.12	0.61
2	5.42	1.56	60.28	29.95	2.79

**Figure S1.** (a) Cross-sectional TEM image of a MoS<sup>2</sup> encapsulated by Al<sub>2</sub>O<sub>3</sub>. Inset: FFT patterns of the MoS<sup>2</sup> obtained in the area marked with the white dashed line. (b) HAADF image of a MoS<sup>2</sup> TFT with Al<sub>2</sub>O<sub>3</sub> passivation; (c) atomic percentage of the atoms contained in the area marked with red and green rectangles in (b).



Figure S2. (a) Raman and (b) XPS spectra of (i) pristine and (ii) UV-ozone treated MoS<sub>2</sub>.

Variations of the physical properties in multilayer MoS<sub>2</sub> with respect to the UV-ozone treatment were investigated using Raman and X-ray photoelectron spectroscopy (XPS). In Figure S2a, both of the two-spectra show two representative peaks belonging to MoS<sub>2</sub>, which are assigned to the E<sup>1</sup><sub>2g</sub> peak at 384 cm<sup>-1</sup> and the A<sub>1g</sub> peak at 409 cm<sup>-1</sup>, respectively [S1\_3]. The Raman spectra exhibit no apparent variation in the peak intensity and position, indicating that negligible defects or lattice disorders are induced by the UV-ozone treatment.

As shown in the XPS spectra (Figure S2b), two peaks with relatively high intensities are observed at 229.3 eV and 232.4 eV for the pristine  $MoS_2$  (i.e., before UV-ozone treatment), which correspond to  $Mo^{4+}$  3d. Another-Other weak peaks corresponding to  $S^{2-}$  2p are identified at 162.1 eV and 163.3 eV [S4,5]. However, after UV-ozone treatment, the intensities of the XPS peaks originatinged from  $Mo^{4+}$  3d and  $S^{2-}$  2p were clearly reduced. In addition, a-strong doublet peaks are observed at the higher binding energies of 232.7 and 235.8 eV<sub>2</sub> implying formation of Mo\_O bonding ( $Mo^{6+}$  3d) [S6]. New peaks also appear at 168.5 and 169.7 eV<sub>2</sub> corresponding to S<sup>6+</sup> 2p (S\_O bonding) [S3]. These results indicate that the  $MoS_2$  surface becomes oxidized during UV-ozone treatment.



**Figure S3.** Transfer characteristics of 14 MoS<sub>2</sub> TFTs before (black line) and after (red line) Al<sub>2</sub>O<sub>3</sub> passivation.



**Figure S4.** Transfer characteristics of six MoS<sub>2</sub> TFTs without Al<sub>2</sub>O<sub>3</sub> passivation before (i.e., pristine) and after UV-ozone treatment.