Supplemental Information for "Combinatorial Insights into Doping Control and Transport Properties of Zinc Tin Nitride"

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Fig. S1 shows conductivity as a function of growth temperature for Zn-rich films with cation composition of 0.60 Zn/(Zn+Sn). Conductivity of Zn-rich films varied by more than an order of magnitude as a function of increasing growth temperature; from 0.1-5.5 S cm⁻¹. This variation of conductivity with temperature was not as pronounced as for variation with composition (Fig. 2a) for a fixed growth temperature. The critical temperature at which the conductivity trend switched from decreasing to increasing was found to be 120°C.

Fig. S2 shows a transmission electron diffraction (TED) pattern from a nominally stoichiometric film grown at 300°C with the six smallest diffraction rings numbered in white (left panel), and a corresponding table giving the calculated lattice spacing from each ring (right panel). The experimental lattice spacing values were found to be consistent with a wurtzite lattice. These values are also consistent with lattice parameters calculated previously for cation-disordered ZnSnN₂, in which the c-axis lattice parameter was 5.53 Å and the a-axis parameter was $3.38 \text{ Å}.^1$

Fig. S3 shows the results of temperature-dependent Hall effect performed on a stoichiometric (Zn/(Zn+Sn) = 0.50) and zinc-rich (Zn/(Zn+Sn) = 0.60) film grown at 230°C, respectively. No significant temperature dependence was observed in the carrier density or

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mobility shown in panel (a), and neither for conductivity shown in panel (b). Considering the relatively low carrier density observed in zinc-rich films, these results support the existence of a large density of competing *n*-type and *p*-type defects combining to yield the ultimate *n*-type carrier density observed.



Fig. S1: Conductivity as determined by 4-point probe paired with thickness measurements, shown as a function of growth temperature for Zn-rich films with 0.60 Zn/(Zn+Sn). Conductivity decreases until 120°C, after which the conductivity increases again. The overall change in conductivity was more than an order of magnitude.

11 Janes State	Diffraction Ring	Lattice Spacing (Å)
	1	2.97
3	2	2.77
	3	2.58
4 2	4	2.02
6 5	5	1.70
	6	1.56

Fig. S2: TED pattern taken from a nominally stoichiometric film grown at 300° C (left) with the six smallest diffraction rings numbered in white. A table giving the corresponding lattice spacing for each diffraction ring is shown on the right. These lattice spacing values are consistent with a wurtzite lattice for ZnSnN₂.



Fig. S3: Temperature-dependent Hall effect data taken from a stoichiometric (0.50) and zinc-rich (0.60) Zn-Sn-N film, respectively. No significant temperature dependence was observed in the carrier density or mobility (a), and neither for conductivity (b). These results support the existence of a large density of competing *n*-type and *p*-type defects combining to yield the ultimate *n*-type carrier density

REFERENCES

Feldberg, N. *et al.* Growth, disorder, and physical properties of ZnSnN₂. *Applied Physics Letters* 103 (2013).