

Table S1: Refinement and lattice parameters of K_2NaAlF_6 phosphor.

Parameters	K_2NaAlF_6 phosphor
Crystal system	Cubic
Space Group	Fm-3m
V	535.11 \AA^3
Lattice Parameters	a = 8.1154 \AA b = 8.1154 \AA c = 8.1154 \AA $\alpha = 90.00^\circ$ $\beta = 90.00^\circ$ $\gamma = 90.00^\circ$
R_{exp}	12.83
R_p	10.22
R_{wp}	17.51
χ^2	2.82



Fig. S1: Surface morphology of the coated solar cell with Eu^{3+} - Tb^{3+} activated/ co-activated K_2NaAlF_6 phosphor has been synthesized by low temperature wet chemical method.

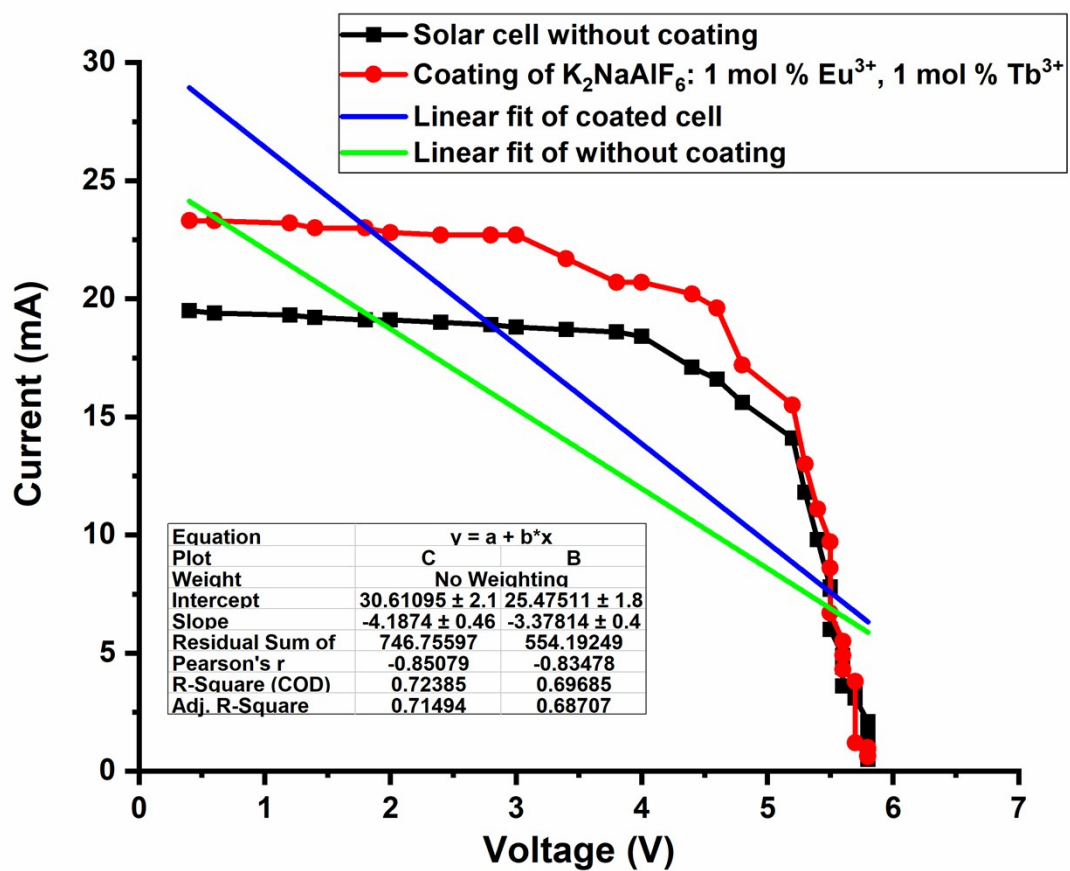


Fig. S2: IV characteristics of Eu^{3+} - Tb^{3+} co-activated K_2NaAlF_6 phosphor under solar simulator.

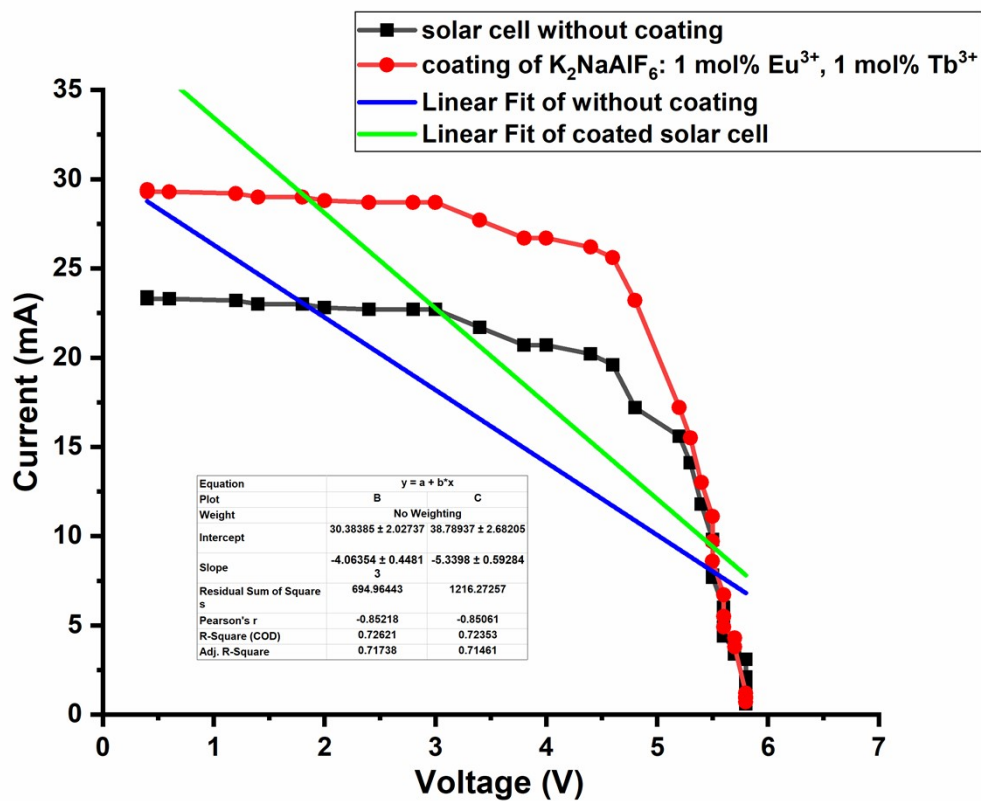


Fig. S3: IV characteristics of Eu^{3+} - Tb^{3+} co-activated K_2NaAlF_6 phosphor under direct sunlight.