Supporting Information

Sodium niobate based hierarchical 3D perovskite nanoparticle clusters

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Equation S1: Scherrer equation for the crystallite size analysis. The analysis of the XRD profile for the evaluation of the average crystallites diameter d_{xrd} was conducted according to the Scherrer equation:

 $d = \frac{0.9\lambda}{FWHMcos\theta}$

The FWHM was evaluated on the (002) diffraction peak.

					Lattice	Reliability
Atom	X	У	Z	Wyck	constants (Å)	factors
Nb	0.75	0.749 (2)	0.751 (1)	4k	a = 5.522 (1)	$R_{P} = 7.02$
Nal	0.75	0	0.247 (6)	2e	b = 7.804(3)	$R_{wp} = 9.38$
Na2	0.75	0.5	0.223 (7)	2f	c = 5.522(1)	$R_{ex} = 4.73$
01	0.25	0	0.128 (5)	2e		
02	0.25	0.5	0.180 (7)	2f		
03	0	0.218 (4)	0.5	4h		
O4	0.5	0.292 (3)	0	4g		

 Table S1: NN300 Pmma (space group number: 51) Refined parameters.



Figure S1: Particle size distribution sample NN300 obtained by SEM microscopy, mean particle size = 312 ± 25 nm.



Figure S2. TEM images and SAED analyses of sample NN270, amorphous nanoparticle cluster (left) and high resolution image of a crystalline mesocrystal (right).



Figure S3. Particle size distribution of the PNPs for sample NN280 (5 \pm 1 nm), NN290 (5 \pm 1 nm) and NN300 (18 \pm 3 nm).



Figure S4: TEM image of sample **NN280**, FFT and profile analysis of the line scan across the fringes. The analyses show the periodic fringe spacing around 0.39 nm, corresponding to the interplanar spacing along the [101] direction of *Pmma* NN.



Figure S5: TEM images of NN300 (scale bar 20 nm); inset: FFT analysis.



Figure S6: Raman spectra evidencing the organic groups in the NN270-300 samples.



Figure S7: NN300 XRD data and Rietveld fitting using the *Pmma* phase.



Figure S8. representation of NN crystal structure along the [010] zone axis



Figure S9. SEM-EDS analysis of NN300 sample, the investigated area is reported in the inset.