

Supplementary information for the article:

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Supplementary file

The low-temperature sonochemical synthesis of up-converting β NaYF₄:Yb,Er mesocrystals

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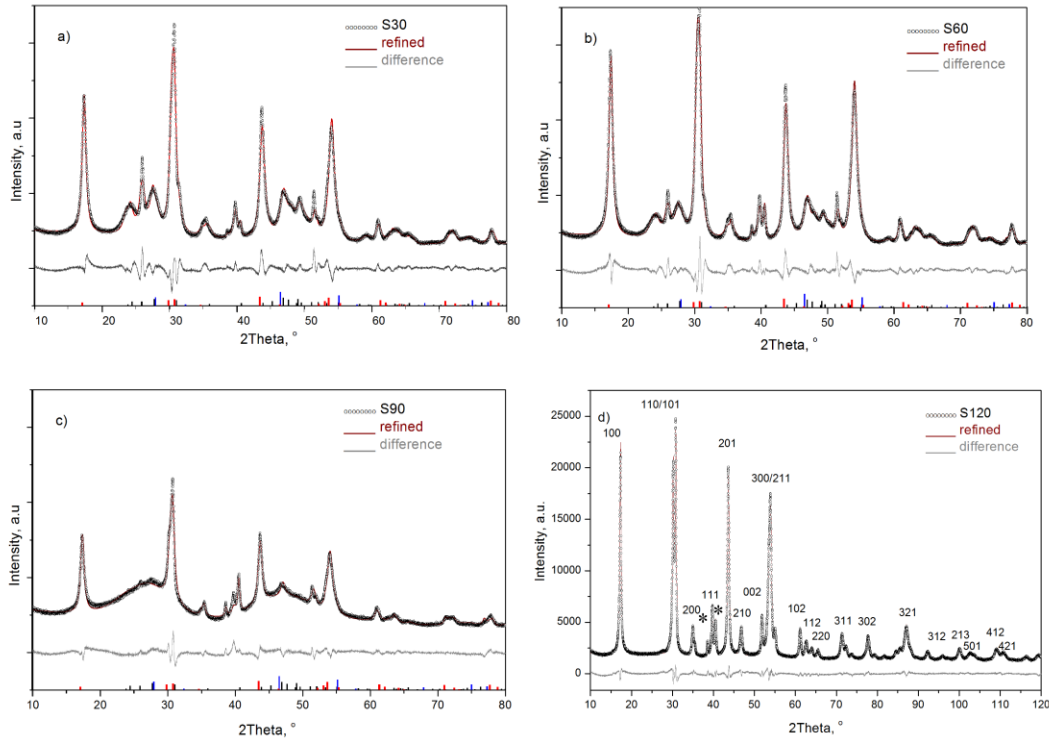


Figure S1. XRPD data and refined patterns of the obtained samples. The Bragg positions of the orthorhombic YF₃ (JCPDS 70-1935) cubic NaYF₄ (JCPDS 77-2042) and hexagonal NaYF₄ (JCPDS 16-0334) are presented as bottom bar lines (orthorhombic-black, cubic-blue, hexagonal-red). Two reflections marked with * in Fig S1d originate from Ti contamination (titanium horn is used during synthesis)

Table 1. Microstructural parameters obtained through Rietveld refinement of Yb,Er doped β -NaYF₄ phase

<i>S.G.</i>	<i>P6₃/m</i>		
<i>a</i> , Å	5.934 (1)		
<i>c</i> , Å	3.529 (1)		
<i>V</i> , Å ³	107.61		
CS, nm	37.6 (4)		
MS	0.31 (1)		
<i>R</i> _{wp}	5.1		
<i>R</i> _{brag}	1.1		
Atom position	x	y	c
Na1/Y1	2/3	2/3	1/4
Na2/Y2	0	0	0.619 (1)
F	0.3098 (3)	0.4009 (4)	1/4

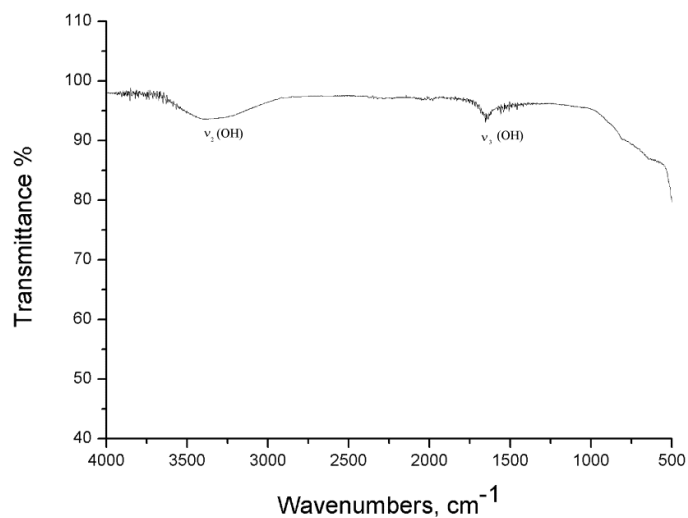


Figure S2. Typical spectrum of FT-IR analysis

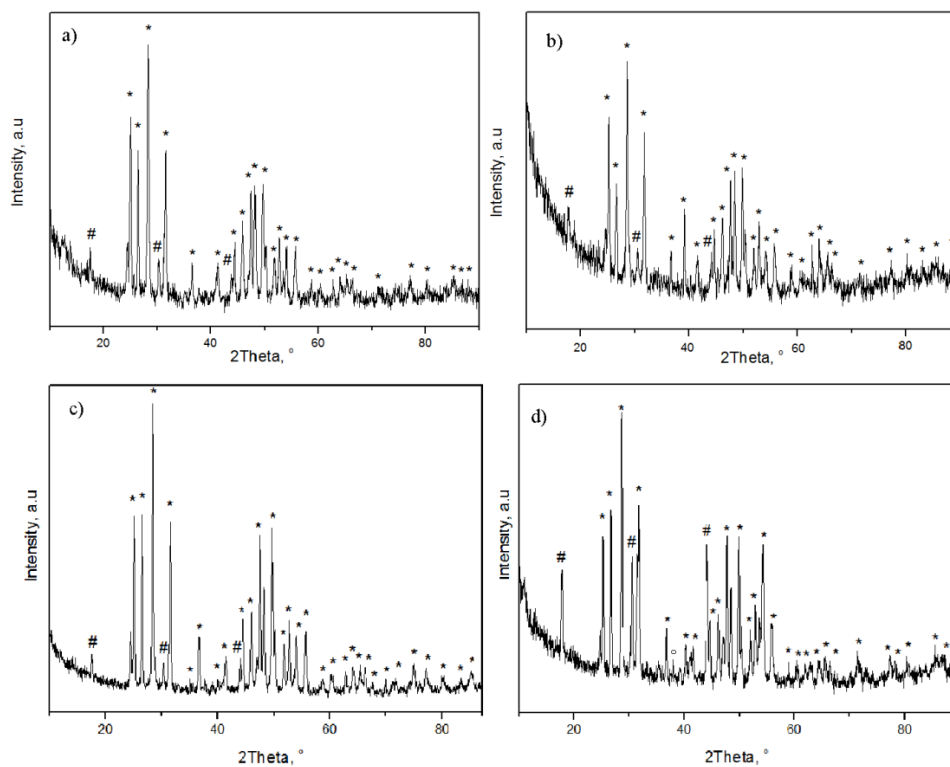


Figure S3. XRPD of samples obtained through sonochemical synthesis after DSC analysis a) S30, b) S60, c) S90, d) S120; XRPD reflections of β - NaYF_4 :Yb,Er are marked with #, while these of YF_3 :Yb,Er with *. NaF presence is indictable in the thermally treated S120 sample due appearance of its strongest reflection at $2\theta=38.85$ marked with $^\circ$ in Fig. S3d.

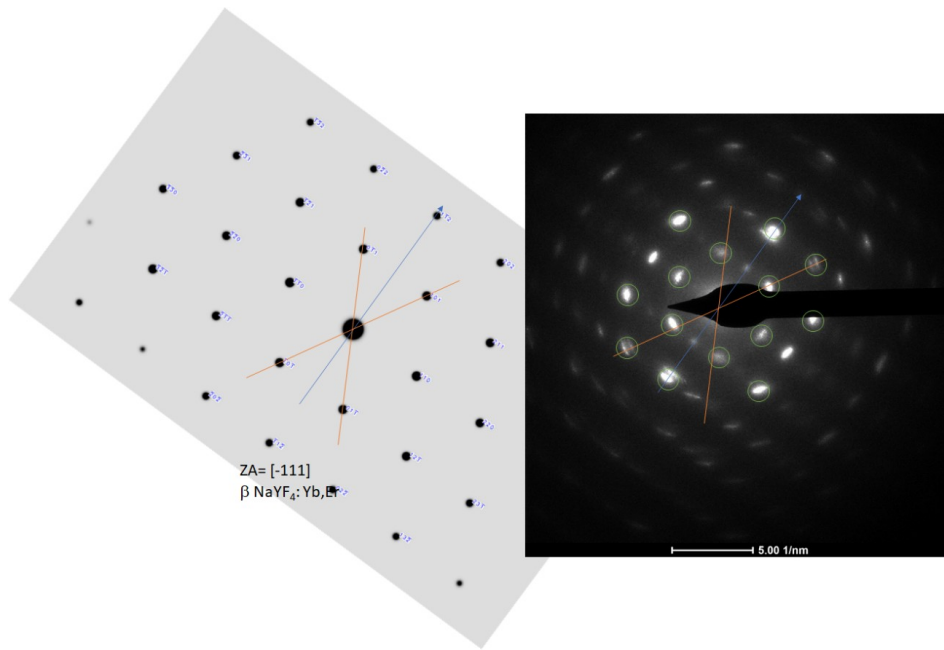


Figure S4: Simulated electron diffraction pattern of β NaYF₄:Yb,Er hexagonal [-111] zone axis compared with experimental SAED. Blue arrow corresponds to growth direction.

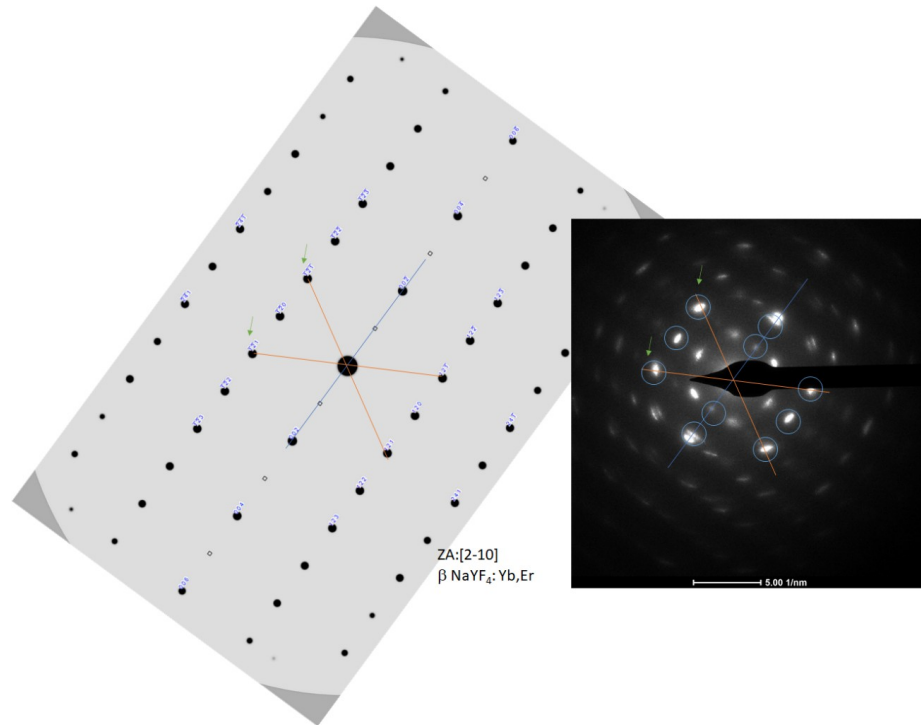


Figure S5: Simulated electron diffraction pattern of β NaYF₄:Yb,Er hexagonal [2-10] zone axis compared with experimental SAED. Note that the (121) planes are present in both zone axis (marked with green arrows). Blue arrow corresponds to growth direction.