

Fig. S1: (a)Experimental compared to linear least-squares fittings of $k^3\chi(k)$ EXAFS oscillations and (b)experimental compared to fitted first-shells filtered EXAFS spectra for unmodified YAG:Tb(20%) powder sintered for 4 h at 1100°C.

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Sample (Heating temperature)		Absorbing atom	N	d (Å)	Sigma ² *10 ³ (Å ²)	R _F (%)
YAG :Tb(20%) R _C =0	xerogel	О	3.5±0.5 (4) 3.1±0.5 (4)	2.33 (2.30) 2.46 (2.43)	4.99 4.99	0.748 e _{not} =1.43
		Al	1.6±0.7 (nd)	3.30 (nd)	16.9	
		Y	0.7±0.4 (3.2)	3.79 (3.67)	8.77	
	200°C	Ο	3.9±0.5 (4) 2.3±0.7 (4)	2.34 (2.30) 2.47 (2.43)	5.60 5.60	1.02 e _{not} =1.47
		Al	1.6±0.8 (nd)	3.29 (nd)	15.41	
		Y	0.6±0.5 (3.2)	3.81 (3.67)	7.95	
	400°C	О	4.0±0.4 (4) 2.3±0.6 (4)	2.32 (2.30) 2.45 (2.43)	9.54 9.54	0.357 e _{not} =1.00
		Al	1.3±0.4 (nd)	3.15 (nd)	14.2	
		Y	1.7±0.5 (3.2)	3.71 (3.67)	13.6	
		Tb	$0.8 \pm 0.6 \ (0.8)$	3.71 (3.67)	13.6	
	600°C	О	3.9±0.4 (4) 1.9±0.6 (4)	2.32 (2.30) 2.45 (2.43)	10.6 10.6	0.283 e _{not} =0.90
		Al	1.4± 0.4 tetra (2)	3.14 (3.00)	11.9	
		Y	1.3±0.9 (3.2)	3.71 (3.67)	12.1	
		Тb	$0.6 \pm 0.9 \ (0.8)$	3.71 (3.67)	12.1	
	700°C	О	4.1±0.4 (4) 2.3±0.6 (4)	2.32 (2.30) 2.44 (2.43)	10.6 10.6	0.329 e _{not} =0.92
		Al	1.3±0.3 tetra (2)	3.14 (3.00)	11.3	
		Y	2.0±0.4 (3.2)	3.71 (3.67)	11.9	
		ТЪ	$0.8 \pm 0.6 \ (0.8)$	3.71 (3.67)	11.9	
	800°C	О	3.0±0.3 (4) 2.8±0.3 (4)	2.31 (2.30) 2.44 (2.43)	4.73 4.73	1.26 e _{not} =1.56
		Al	1.2± 0.3 tetra (2) 2.1±1.0 octa (4) 2.1±0.8 (4)	3.04 (3.00) 3.40 (3.35) 3.71 (3.67)	4.22 11.6 4.88	
		Y	2.8±0.3 (3.2)	3.71 (3.67)	4.88	
		Тb	1.2±0.3 (0.8)	3.71 (3.67)	4.88	
	1100°C	0	4 (4) 4 (4)	2.32 (2.30) 2.46 (2.43)	5.70 5.70	3.56 e _{not} =1.83
		Al	2 tetra (2) 4 octa (4) 4 (4)	3.04 (3.00) 3.38 (3.35) 3.71 (3.67)	3.24 9.26 2.66	
		Y	3 (3.2)	3.70 (3.67)	2.66	
		Tb	1 (0.8)	3.70 (3.67)	2.66	

Table S1: Structural parameters determined from the EXAFS data recorded at Tb edge for unmodified samples.

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Sample (Heating temperature)		Absorbing atom	N	d (Å)	Sigma ² *10 ³ (Å ²)	R _F (%)
	xerogel	О	5.0±0.5 (4) 2.6±0.9 (4)	2.34 (2.30) 2.47 (2.43)	6.48 6.48	1.03 e _{not} =0.93
		Al	1.0±0.6 (nd)	3.33 (nd)	13.3	
		Y	0.4±0.5 (3.2)	3.82 (3.67)	6.73	
	200°C	Ο	4.2±0.5 (4) 2.9±0.7 (4)	2.34 (2.30) 2.47 (2.43)	6.40 6.40	0.768 e _{not} =1.30
		Al	1.1±0.6 (nd)	3.27 (nd)	12.0	
		Y	0.8±0.4 (3.2)	3.77 (3.67)	7.07	
		Tb	0.6±0.6 (3.2)	3.77 (3.67)	7.07	
	400°C	Ο	3.2±0.8 (4) 2.8±1.0 (4)	2.31(2.30) 2.44 (2.43)	6.33 6.33	0.678 e _{not} =1.09
		Al	0.9±0.2 (nd)	3.20 (nd)	10.4	
		Y	0.6±0.5 (3.2)	3.71 (3.67)	7.27	
YAG :Tb(20%) R _C =1		Tb	0.2±0.8 (0.8)	3.71 (3.67)	7.27	
	600°C	О	2.6±0.5 (4) 2.2±0.6 (4)	2.29 (2.30) 2.42 (2.43)	5.80 5.80	1.05 e _{not} =1.18
		Al	0.9±0.2 tetra (2)	3.11 (3.00)	5.22	
		Y	1.4±0.3 (0.3)	3.69 (3.67)	5.74	
		Tb	0.9±0.4 (0.8)	3.69 (3.67)	5.74	
	700°C	Ο	2.5±0.5 (4) 2.5±0.5 (4)	2.29 (2.30) 2.42 (2.43)	4.72 4.72	0.580 e _{not} =1.10
		Al	1.1±0.2 tetra (2)	3.06 (3.00)	6.32	
		Y	2.8±0.3 (3.2)	3.71 (3.67)	5.70	
		Tb	1.4±0.4 (0.8)	3.71 (3.67)	5.70	
	800°C	Ο	2.8±0.6 (4) 3.6±0.9 (4)	2.31 (2.30) 2.44 (2.43)	3.84 3.84	2.21 e _{not} =1.80
		Al	2.0±0.3 tetra (2) 4.1±0.9 octa (4) 3.9±0.3 (4)	3.04 (3.00) 3.40 (3.35) 3.71 (3.67)	3.98 9.58 3.84	
		Y	2.7±0.6 (3.2)	3.71 (3.67)	3.84	
		Tb	1.3±0.6 (0.8)	3.71 (3.67)	3.84	
	1100°C	Ο	4 (4) 4 (4)	2.32 (2.30) 2.45 (2.43)	6.2 6.2	2.94 e _{not} =1.80
		Al	2 tetra (2) 4 octa (4) 4 (4)	3.04 (3.00) 3.38 (3.35) 3.71 (3.67)	2.58 8.01 3.55	
		Y	3 (3.2)	3.70 (3.67)	3.55	
		Tb	1 (0.8)	3.70 (3.67)	3.55	

Table S2: Structural parameters determined from the EXAFS data recorded at Tb edge for modified samples.



Fig. S2: PRDFs uncorrected from phase shift for TAG and acac-modified YAG powders(R_C =1) sintered at different temperatures: X – xerogel, B – 200°C, K – 400°C, I – 600°C, H – 700°C, C – 800°C and A – 1100°C.