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Electronic Supplementary Information

SynthesisofPerovskite-typeManganites Yb_{1-} $_xDy_xMnO_3$ (0.1 $\leq x \leq 0.5$)viaSolid-StateReactionandHigh-pressureFluxMethodFollowedby,StructuralCharacterizationandMagnetic

Property Studies

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Yb(1)-O(1)#1	2.236(4)	O(1)#1-Yb(1)-O(2)	100.87(10)
Yb(1)-O(2)	2.279(3)	O(1)#1-Yb(1)-O(2)#2	100.87(10)
Yb(1)-O(2)#2	2.279(3)	O(2)-Yb(1)-O(2)#2	78.17(13)
Yb(1)-O(1)	2.304(4)	O(1)#1-Yb(1)-O(1)	91.29(9)
Yb(1)-O(2)#3	2.515(3)	O(2)-Yb(1)-O(1)	138.46(7)
Yb(1)-O(2)#4	2.515(3)	O(2)#2-Yb(1)-O(1)	138.46(7)
Yb(1)-O(2)#5	2.562(3)	O(1)#1-Yb(1)-O(2)#3	137.99(8)
Yb(1)-O(2)#6	2.562(3)	O(2)-Yb(1)-O(2)#3	118.85(5)
Mn(2)-O(2)#10	1.908(3)	O(2)#2-Yb(1)-O(2)#3	76.24(6)
Mn(2)-O(2)#4	1.908(3)	O(1)-Yb(1)-O(2)#3	68.58(10)
Mn(2)-O(1)	1.9467(14)	O(1)#1-Yb(1)-O(2)#4	137.99(8)
Mn(2)-O(1)#11	1.9467(14)	O(2)-Yb(1)-O(2)#4	76.24(6)
Mn(2)-O(2)#6	2.227(3)	O(2)#2-Yb(1)-O(2)#4	118.85(5)
Mn(2)-O(2)#12	2.227(3)	O(1)-Yb(1)-O(2)#4	68.58(10)
		O(2)#3-Yb(1)-O(2)#4	69.69(12)
		O(1)#1-Yb(1)-O(2)#5	69.01(7)
		O(2)-Yb(1)-O(2)#5	150.11(6)

Table S1. Bond lengths [Å] and angles [deg] for the *o*-Yb_{0.5}Dy_{0.5}MnO₃

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O(2)#2-Yb(1)-O(2)#5	76.47(11)
O(1)-Yb(1)-O(2)#5	71.19(7)
O(2)#3-Yb(1)-O(2)#5	69.64(4)
O(2)#4-Yb(1)-O(2)#5	130.52(5)
O(1)#1-Yb(1)-O(2)#6	69.01(7)
O(2)-Yb(1)-O(2)#6	76.47(11)
O(2)#2-Yb(1)-O(2)#6	150.11(6)
O(1)-Yb(1)-O(2)#6	71.19(7)
O(2)#3-Yb(1)-O(2)#6	130.52(5)
O(2)#4-Yb(1)-O(2)#6	69.64(4)
O(2)#5-Yb(1)-O(2)#6	121.65(13)
O(2)#10-Mn(2)-O(2)#4	180.00(15)
O(2)#10-Mn(2)-O(1)	90.23(14)
O(2)#4-Mn(2)-O(1)	89.77(14)
O(2)#10-Mn(2)-O(1)#11	89.77(14)
O(2)#4-Mn(2)-O(1)#11	90.23(14)
Mn(2)-O(1)- Mn(2)	141.3(2)
O(2)#10-Mn(2)-O(2)#6	91.30(5)
O(2)#4-Mn(2)-O(2)#6	88.70(5)
O(1)-Mn(2)-O(2)#6	85.48(13)
O(1)#11-Mn(2)-O(2)#6	94.52(13)
O(2)#10-Mn(2)-O(2)#12	88.70(5)
O(2)#4-Mn(2)-O(2)#12	91.30(5)
O(1)-Mn(2)-O(2)#12	94.52(13)
O(1)#11-Mn(2)-O(2)#12	85.48(13)
 Mn(2)-O(2)- Mn(2)	143.42(14)

Symmetry transformations used to generate equivalent atoms: #1 x+1/2,y,-z-1/2 #2 x,-y+1/2,z #3 x-1/2,-y+1/2,-z+1/2 #4 x-1/2 y -z+1/2 #5 -x+1 y-1/2 -z #6 -x+1 -y+1 -z

#4 x-1/2,y,-z+1/2	#5	-x+1,y-1/2,-z	#6 ·	-x+1,-y+1,-z
#10 -x+1/2,-y+1,z-1	/2	#11 -x,-y+1,-z		#12 x-1,y,z

Table S2. Atomic coordinates and equivalent isotropic displacementparameters (Ų) for Yb_{0.5}Dy_{0.5}MnO₃

	X	у	Z	U(eq)
Yb(1)	4161(1)	2500	173(1)	5(1)
Dy(1)	4161(1)	2500	173(1)	5(1)
Mn(2)	0	5000	0	3(1)
O(1)	396(6)	2500	-1146(8)	5(1)
O(2)	6707(4)	4455(3)	1999(6)	7(1)