Supplementary Information

Optimization of fluxes for Yb³⁺:YMgB₅O₁₀ crystal growth and intense multiwavelength emission characteristics in spectra and laser performances

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Figure S1. Schematic diagram of crystal growth apparatus.

Figure S2. Thermal expansion coefficients of Yb:YMB crystal.

Figure S3. Differential thermal analysis (DTA) and thermogravimetry (TG) of Yb:YMB crystal.

Figure S4. Powder XRD patterns of Yb:YMB crystal.

Figure S5. Cyan light emission phenomenon in the laser experiment.

Figure S6. Laser spectra corresponding to the cyan light emission.

 Table S1 Atomic coordinates and equivalent isotropic displacement parameters of

 Yb:YMB crystal.

Table S2 Selected bond distances (Å) of Yb:YMB crystal.



Fig. S1. Schematic diagram of crystal growth apparatus. (1) seed-holder, (2) watch window, (3) furnace cover (thermal insulation materials), (4) nickel-chrome heating wires, (5) Al_2O_3 tube, (6) platinum wires, (7) seed, (8) platinum crucible, (9) thermocouple, (10) solution, (11) thermal insulation materials (the height is based on the thermal field), (12) thermal insulation materials.



Fig. S2. Thermal expansion coefficients of Yb:YMB crystal.



Fig. S3. Differential thermal analysis (DTA) and thermogravimetry (TG) of Yb:YMB crystal.



Fig. S4. Powder XRD patterns of Yb:YMB crystal.



Fig. S5. Cyan light emission phenomenon in the laser experiment.



Fig. S6. Laser spectra corresponding to the cyan light emission.

Atom	x/a	<i>y</i> /b	z/c	$U(eq)[Å^2]^*$
Y(1)	1855(1)	1866(1)	7395(1)	4(1)
Yb(1)	1855(1)	1866(1)	7395(1)	4(1)
Mg(1)	6002(1)	4080(2)	6280(1)	6(1)
B(1)	-653(4)	3113(5)	9105(4)	5(1)
B(2)	5842(4)	751(5)	7430(4)	5(1)
B(3)	-2234(4)	326(5)	5548(4)	5(1)
B(4)	3398(4)	-1003(5)	9939(4)	6(1)
B(5)	-177(4)	1740(5)	3955(4)	3(1)
O(1)	1767(2)	-1075(3)	5822(2)	5(1)
O(2)	456(2)	2710(3)	5210(2)	5(1)
O(3)	-94(2)	4091(3)	8024(2)	6(1)
O(4)	58(2)	1504(3)	9262(2)	5(1)
O(5)	6860(2)	327(3)	8741(2)	5(1)
O(6)	2569(2)	-517(3)	8735(2)	6(1)
O(7)	6816(2)	1509(3)	6353(2)	5(1)
O(8)	-841(2)	-281(3)	6484(2)	5(1)
O(9)	3184(2)	1263(3)	4910(2)	5(1)
O(10)	4636(2)	2082(3)	7700(2)	4(1)

Table S1 Atomic coordinates and equivalent isotropic displacement parameters of Yb:YMB crystal

U(eq) is defined as one-third of the trace of the orthogonalized U_{ij} tensor.

Table S2 Selected bond distant	nces (Å) of Yb:YMB crystal
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Y(1)/Yb(1)-O(6)	2.255(2)	B(1)-O(9)#4	1.365(4)
Y(1)/Yb(1)-O(6)#1	2.311(2)	B(2)-O(3)#5	1.458(4)
Y(1)/Yb(1)-O(10)	2.373(2)	B(2)-O(7)	1.462(4)
Y(1)/Yb(1)-O(2)	2.384(2)	B(2)-O(10)	1.471(4)
Y(1)/Yb(1)-O(4)	2.409(2)	B(2)-O(5)	1.490(4)
Y(1)/Yb(1)-O(3)	2.461(2)	B(3)-O(7)#6	1.449(4)
Y(1)/Yb(1)-O(1)#1	2.516(2)	B(3)-O(1)#7	1.478(4)
Y(1)/Yb(1)-O(1)	2.664(2)	B(3)-O(9)# 7	1.496(4)
Y(1)/Yb(1)-O(9)	2.687(2)	B(3)-O(8)	1.500(4)
Mg(1)-O(4)#2	2.050(2)	B(4)-O(6)	1.341(4)
Mg(1)-O(5)#3	2.051(2)	B(4)-O(5)#8	1.366(4)
Mg(1)-O(7)	2.063(3)	B(4)-O(2)# 5	1.391(4)
Mg(1)-O(4)#1	2.091(2)	B(5)-O(2)	1.457(4)
Mg(1)-O(8)#1	2.158(2)	B(5)-O(1)#7	1.471(4)
Mg(1)-O(10)	2.365(2)	B(5)-O(10)#9	1.473(4)
B(1)-O(3)	1.362(4)	B(5)-O(8)#7	1.478(4)
B(1)-O(4)	1.363(4)		

Symmetry codes: (#1) -x+1/2,y+1/2,-z+3/2; (#2) x+1/2,-y+1/2,-z-1/2; (#3) -x+3/2,y+1/2,-z+3/2; (#4) x-1/2,-y+1/2,z+1/2; (#5) -x+1/2,y-1/2,-z+3/2; (#6) x-1,y,z; (#7) -x,-y,-z+1; (#8) -x+1,-y,-z+2; (#9) x-1/2,-y+1/2,z-1/2