

Additional Materials

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>Lc1|C. unicolor|383410|pI: 5.21|Mw: 55.9 kDa
MSFNAAFATFLTLAALASVYAALGPVTDLHIVNT35NISTDGFTRPAVLAGGTFPGPVITAKKGNFQVNVIDDLTNEAMLKSTSIHWHGLFQKGTNWADGP      100
AFVNSQCPIISGNSFLYEFVSPDQAGTYWYHSHLSTQYCDGLRGLVLIYDPPDPHAGLYDVDDENTIIITLADWYHTFARQIPGLPIADTTILINGLGRNH199NG      200
SADAELSVINVHAGKRYRFRFLVSI238CDPNYVFSID238DMTVIEVEGTVNPKPLTVDSIQIFAAQRYSFVLHADRPAGNYWIRANPSFGNTGFAGHINSAIL      300
RYDGAPCPEPHTRQTNSTKFLNEIDLHPLVPIPVPGKPYGGVDVVKNLVFNFTDGHYTI361INNATFTPPSVVLLQILSGTTS361SAQGLLP361SGDVIELPLGKT      400
VEITLAAGVIGGPHFHLHGHSFHVVRSACQNTSNYIDPIVRDVTNIGMGDNVIRFTTDNPGPWFLHCHIDWHLEAGFAVVLAEGMNETRAANPTPDE      500
WDSLCDTYNALPSCDQ*      516

>Lc2|C. unicolor|408157|pI: 5.13|Mw: 57.6 kDa
MGSKSSILTFITFALALGSYAAIGPVADLHIVNKDLAPDGVTRPILAGGTFPGPLITGQKGNFKLNVIDELTDERMLTPTS35SIHWHGFFQKGTNWADGP      100
AFVNSQCPIIPDNSFLYDFNVPDQAGTFWYHSHLSTQYCDGLRGAFVVYDPEDPHKDLYDVDDSTVITLADWYHVLAPT199VKFTATPDSTLINGLGRSHDG      200
PADAELAVIEVEQ238GKRYRFRFLVNI238GCFPNYDFSID238GMITIEVEGVNSQ361PHTVDSIQIFVAQRYSFVLNANQ361PVDNYVVKAVPNLGDKST361DGGINSAIL      300
RYKGA361PVAEPKTRNVKSRM361PLETDLHPLERMPVPGREVPVGGADIVHLLDL361SFSSDRFFI361NGSSYRD361PPVVLLQILSGVKT361AQDLLPAGNVIGLELGKV      400
VEV361IIPRAHDGPHFHLHGHNFWVRSAGSKEYNFDDPILRDVVSIGGGNDEV361IRFVTDNAGPWFLHCHIDWHLEAGLAVVFAEG361IESTAAANPTPQQ      500
WDELCPKYNALIQSQRALINNSTVN*      525
    
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Figure S1 Protein sequences of Lc1 and Lc2 from *Cerrena unicolor*. Amino acids in bold indicate a putative signal peptide. Predicted N-glycosylation sites are in blue.

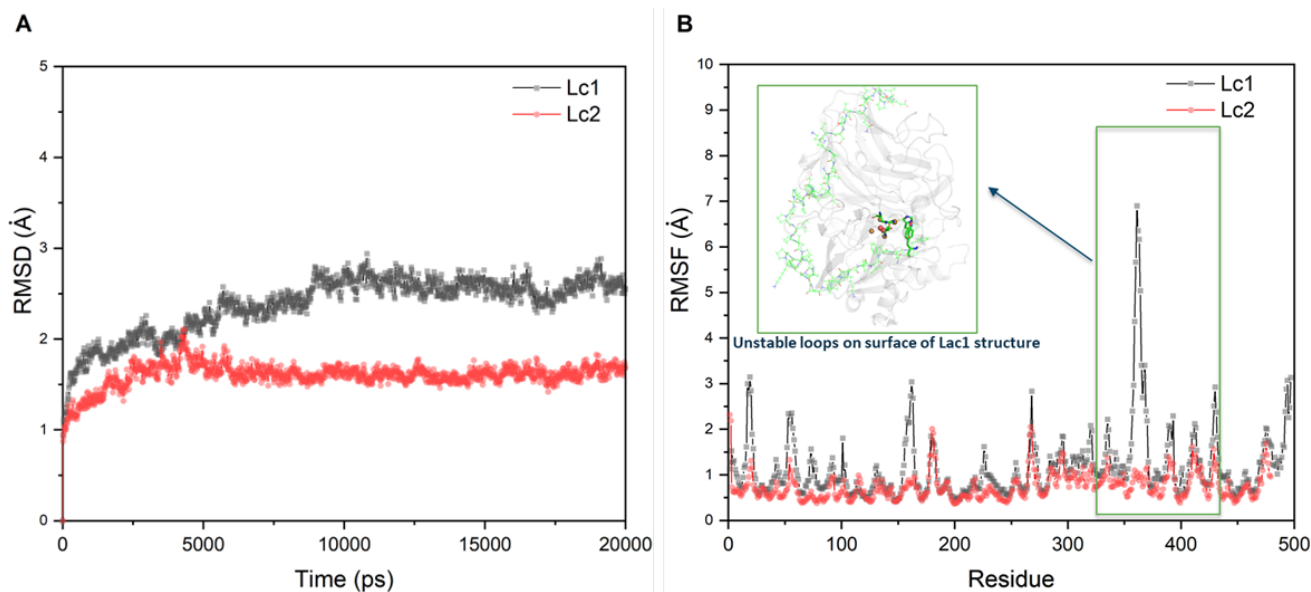


Figure S2 Time course of RMSD for Lc1 and Lc2 (A) and per-residue RMSF flexibility (B) of Lc1 and Lc2 for 50 ns of MD simulation at 333.15 K. The inset image shows the surface loop for which the RMSF larger in Lc1.

Table S1 Salt-bridge network analysis for structural models of Lc1 and Lc2. The analysis was conducted using VMD software. Salt bridges were as defined as a negatively charged oxygen atom of an acidic residue being within 3.2 Å of a positively charged nitrogen atom of basic residue.

Lc1: 35 salt-bridge bonds		Lc2: 43 salt-bridge bonds	
Non-homologous salt-bridges	Homologous salt-bridges	Homologous salt-bridges	Non-homologous salt-bridges
Asp ₆₃ - Lys ₆₀	Asp ₁₂₂ - Lys ₉₂	Asp ₁₂₂ - Lys ₉₂	Asp ₃₅ - Arg ₇₇
Asp ₁₂₂ - Lys ₆₁	Asp ₁₃₉ - Arg ₄₃	Asp ₁₃₉ - Arg ₄₃	Asp ₃₅ - Lys ₃₄
Asp ₁₅₁ - Lys ₆₁	Asp ₁₄₉ - Lys ₆₁	Asp ₁₄₉ - Lys ₆₁	Asp ₇₅ - Arg ₇₇
Asp ₁₅₉ - Lys ₂₁₅	Asp ₁₅₂ - Arg ₂₁₈	Asp ₁₅₂ - Arg ₂₁₈	Asp ₁₁₇ - Lys ₆₆
Asp ₂₃₆ - Arg ₂₇₃	Asp ₁₅₉ - Arg ₂₁₆	Asp ₁₅₉ - Arg ₂₁₆	Asp ₁₅₆ - Arg ₂₁₆
Asp ₂₃₆ - Arg ₃₁₃	Asp ₁₆₁ - Arg ₂₁₈	Asp ₁₆₁ - Arg ₂₁₈	Asp ₁₅₆ - Lys ₁₅₅
Asp ₂₇₂ - Arg ₂₇₃	Asp ₁₆₁ - Arg ₂₂₀	Asp ₁₆₁ - Arg ₂₂₀	Asp ₂₀₃ - Arg ₁₉₆
Asp ₃₂₅ - Lys ₃₁₉	Asp ₁₇₁ - Arg ₁₄₂	Asp ₁₇₁ - Arg ₁₄₂	Asp ₂₃₁ - Lys ₂₈₉
Asp ₅₁₅ - Lys ₈₀	Asp ₂₃₅ - Arg ₂₈₁	Asp ₂₃₅ - Lys ₂₈₁	Asp ₂₃₁ - Lys ₃₁₆
Glu ₇₆ - Lys ₈₀	Asp ₂₅₅ - Lys ₃₁₉	Asp ₂₅₅ - Arg ₃₁₈	Asp ₂₇₆ - Lys ₃₁₁
Glu ₂₄₅ - Arg ₂₆₄	Asp ₃₄₄ - Lys ₃₃₇	Asp ₃₄₄ - Arg ₃₃₇	Asp ₂₉₂ - Arg ₃₁₃
	Asp ₃₄₄ - Lys ₃₉₉	Asp ₃₄₄ - Lys ₃₉₉	Asp ₃₅₀ - Arg ₄₀₈
	Asp ₄₃₈ - Lys ₂₅₀	Asp ₄₃₈ - Arg ₃₃₁	Asp ₃₅₆ - Lys ₁₈₁
	Asp ₄₄₃ - Arg ₂₆₄	Asp ₄₄₃ - Arg ₂₆₄	Asp ₃₅ - Arg ₇₇
	Asp ₄₅₂ - Arg ₄₂₇	Asp ₄₅₂ - Arg ₄₂₇	Asp ₃₅ - Lys ₃₄
	Asp ₄₆₁ - Lys ₉₂	Asp ₄₆₁ - Lys ₉₂	Asp ₃₆₇ - Arg ₃₆₆
	Glu ₂₄₃ - Arg ₂₁₈	Glu ₂₄₃ - Arg ₂₁₈	Asp ₄₃₇ - Arg ₃₃₁
	Glu ₂₄₃ - Arg ₂₂₀	Glu ₂₄₃ - Arg ₂₂₀	Asp ₄₅₂ - Arg ₄₀₈
	Glu ₂₄₅ - Arg ₄₄₂	Glu ₂₄₅ - Arg ₄₄₂	Asp ₄₅₂ - Lys ₄₃₂
	Glu ₃₀₉ - Arg ₁₉₆	Glu ₃₀₉ - Arg ₁₉₆	
	Glu ₃₂₃ - Arg ₄₄₂	Glu ₃₂₃ - Arg ₄₄₂	
	Glu ₃₉₄ - Arg ₄₉₂	Glu ₃₉₆ - Lys ₃₉₉	
	Glu ₄₀₂ - Arg ₄₅₇	Glu ₄₀₂ - Arg ₄₅₇	
	Glu ₄₇₇ - Arg ₁₇₈	Glu ₄₇₇ - Lys ₁₈₁	