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Supplemental Information

Structural Basis for RNA Replication

by the SARS-CoV-2 Polymerase

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Number	Constructs	Oligonucleotide sequences			
1	T33-1/P10	GGGAGAUGAAAGUCUCCA <mark>CCUGUG</mark> UCGUCGAAA			
2	T33-2/P10	GGGAGAUGAAAGUCUCCA <mark>GCAUAU</mark> UCGUCGAAA			
3	T33-3/P10	GGGAGAUGAAAGUCUCCA <mark>CCGGG</mark> GUCGUCGAAA			
4	T33-4/P10	GGGAGAUGAAAGUCUCCA <mark>CCGUGU</mark> UCGUCGAAA			
5	T33-5/P10	GGGAGAUGAAAGUCUCCA <mark>CCUUUU</mark> UCGUCGAAA			
6	T33-6/P10	GGGAGAUGAAAGUCUCCA <mark>GGUCUC</mark> UCGUCGAAA			
7	T31/P8	GGGAGAUGAAAGUCUCCA <mark>GGUCUC</mark> GUGGAAA			
8	T34/P10	GGGAGAUGAAAGUCUCCA <mark>GGCUCUC</mark> UCGUCGAAA			
9	T35/P10	GGGAGAUGAAAGUCUCCA <mark>GGUCUCU</mark> CGUCGAAA			
10	T36/P10	GGGAGAUGAAAGUCUCCA <mark>GGCUCUCU</mark> CGUCGAAA			
11	T37/P10	GGGAGAUGAAAGUCUCCA <mark>GGUCUCUCU</mark> CGUCGAAA			
12	T33-7/P10	GGGAGAUGAAAGUCUCCA <mark>GGUCUCUCU</mark> CGUCGAAA			
N/A	P10	UGUUCGACGA			
N/A	P8	UGUUCCGA			

 Table S1. Oligonucleotide sequences information, Related to STAR Methods.

	Pre-trai	Post- translocated catalvtic				
	Overall	Conformation I	Conformation II	complex		
PDB entry	7C2K	N/A	N/A	7BZF		
EMDB entry	EMD-30275	EMD-30283	EMD-30284	EMD-30252		
Data collection and processing						
Magnification	165,000	165,000	165,000	165,000		
Voltage (keV)	300	300	300	300		
Electron exposure (e ⁻ /A ²)	60.00	60.00	60.00	60.00		
Defocus range (µm)	-2.0 to -1.0	-2.0 to -1.0	-2.0 to -1.0	-1.8 to -1.0		
Pixel size (A)	0.82	0.82	0.82	0.82		
Symmetry imposed	C1	C1	C1	C1		
Initial particle images	1,330,896	1,330,896	1,330,896	1,235,162		
(NO.)	102 150	65 462	27 606	110 662		
Map global resolution $(Å)$	2 03	3 03	37,090	3.26		
Clobal resolution ESC	2.95	0.142	0.12	0.142		
threshold	0.143	0.145	0.145	0.145		
Man sharpening <i>B</i> factor	97.6	86 1	76.4	90 0		
$(Å^2)$	57.0	00.1	70.4	55.0		
Map local resolution	1.8-7.5	1.8-10.0	1.8-10.0	2.0-8.0		
range (Å)						
Model resolution (Å)	2.8	N/A	N/A	3.2		
FSC threshold	0.143	N/A	N/A	0.143		
Model resolution range	co to 2.8	N/A	N/A	co to 3.2		
(Å)						
Model composition						
Non-hydrogen atoms	10,692	N/A	N/A	9,932		
Protein residues	1,256	N/A	N/A	1,200		
Nucleotide residues	31	N/A	N/A	19		
Ligands	3	N/A	N/A	2		
<i>B</i> factors (Å ²)						
Protein	45.91	N/A	N/A	51.33		
Nucleotide	78.99	N/A	N/A	74.57		
Ligand	22.30	N/A	N/A	42.29		
R.m.s. deviations						
Bond lengths (A)	0.005	N/A	N/A	0.004		
Bond angles (°)	0.645	N/A	N/A	0.988		
CC(mask)	0.90	N/A	N/A	0.78		
CC(box)	0.79	N/A	N/A	0.73		
Validation						
MolProbity score	1./1	N/A	N/A	2.09		
Clashscore	7.25	N/A	N/A	11.67		
Poor rotamers (%)	0.00	N/A	N/A	0.00		
	05 51	N1/A	N1/A	01 50		
Favored (%)	95.51	N/A	N/A	91.50		
Allowed (%)	4.49	N/A	N/A	8.50		
Disallowed (%)	0.00	N/A	N/A	0.00		
Model coverage	A.4. \/007.	N1/A	N1/A	A.4. T000		
Chain A	A1~V9U/;	N/A	IN/A	AI~I090; T012~T020		
Chain P	0910~1929 976~N100	NI/A	NI/A	1912~1929 K72~N102		
	D0~070	Ν/ <i>Γ</i> λ ΝΙ/Λ	Ν/ <i>Γ</i> λ	S1~160		
	FU~U/Z	N/A	IN/ <i>F</i> A	61~100 K107-A101		
	A00~A191	IN/A	IN/A	NIZ/~A191		