



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:57 AM UTC

PDB ID : 8FMA / pdb_00008fma
EMDB ID : EMD-29290
Title : Nodavirus RNA replication proto-crown, detergent-solublized C11 multimer
Authors : Zhan, H.; Unchwaniwala, N.; Rebolledo Viveros, A.; Pennington, J.; Horswill, M.; Broadberry, R.; Myers, J.; den Boon, J.; Grant, T.; Ahlquist, P.
Deposited on : 2022-12-22
Resolution : 3.10 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

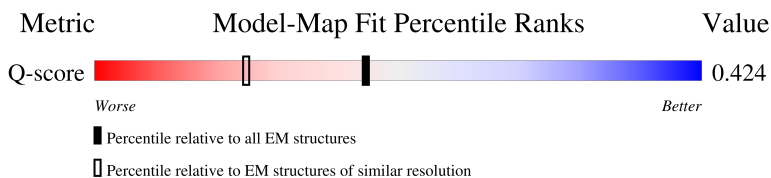
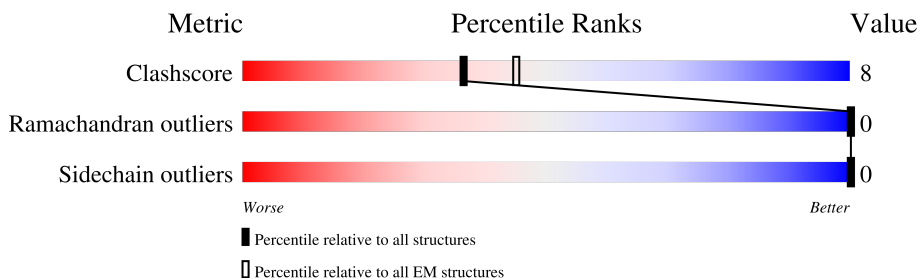
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



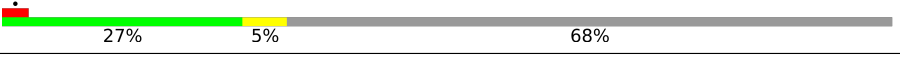
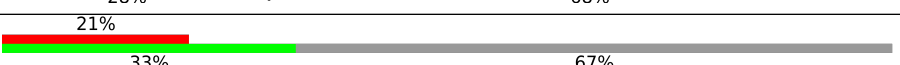
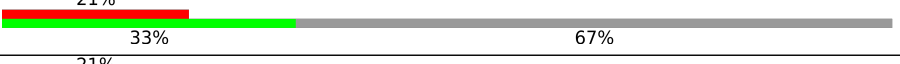
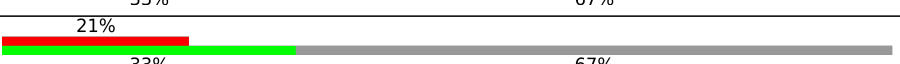
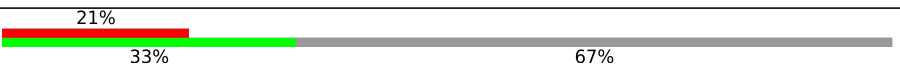
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14724 (2.60 - 3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1011	
1	B	1011	
1	C	1011	
1	D	1011	

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Mol	Chain	Length	Quality of chain
1	E	1011	
1	F	1011	
1	G	1011	
1	H	1011	
1	I	1011	
1	J	1011	
1	K	1011	
1	L	1011	
1	M	1011	
1	N	1011	
1	O	1011	
1	P	1011	
1	Q	1011	
1	R	1011	
1	S	1011	
1	T	1011	
1	U	1011	
1	V	1011	

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 72974 atoms, of which 28633 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	B	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	C	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	D	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	E	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	F	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	G	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	H	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	I	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	J	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	K	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	A	324	Total	C	H	N	O	S	0	0
			5223	1673	2603	459	479	9		
1	L	333	Total	C	N	O			0	0
			1411	724	338	349				
1	M	333	Total	C	N	O			0	0
			1411	724	338	349				
1	N	333	Total	C	N	O			0	0
			1411	724	338	349				
1	O	333	Total	C	N	O			0	0
			1411	724	338	349				
1	P	333	Total	C	N	O			0	0
			1411	724	338	349				
1	Q	333	Total	C	N	O			0	0
			1411	724	338	349				

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Mol	Chain	Residues	Atoms				AltConf	Trace
1	R	333	Total	C	N	O	0	0
			1411	724	338	349		
1	S	333	Total	C	N	O	0	0
			1411	724	338	349		
1	T	333	Total	C	N	O	0	0
			1411	724	338	349		
1	U	333	Total	C	N	O	0	0
			1411	724	338	349		
1	V	333	Total	C	N	O	0	0
			1411	724	338	349		

There are 308 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	897	LEU	MET	conflict	UNP Q66929
B	999	SER	-	expression tag	UNP Q66929
B	1000	ARG	-	expression tag	UNP Q66929
B	1001	LEU	-	expression tag	UNP Q66929
B	1002	GLU	-	expression tag	UNP Q66929
B	1003	GLU	-	expression tag	UNP Q66929
B	1004	GLU	-	expression tag	UNP Q66929
B	1005	LEU	-	expression tag	UNP Q66929
B	1006	ARG	-	expression tag	UNP Q66929
B	1007	ARG	-	expression tag	UNP Q66929
B	1008	ARG	-	expression tag	UNP Q66929
B	1009	LEU	-	expression tag	UNP Q66929
B	1010	THR	-	expression tag	UNP Q66929
B	1011	GLU	-	expression tag	UNP Q66929
C	897	LEU	MET	conflict	UNP Q66929
C	999	SER	-	expression tag	UNP Q66929
C	1000	ARG	-	expression tag	UNP Q66929
C	1001	LEU	-	expression tag	UNP Q66929
C	1002	GLU	-	expression tag	UNP Q66929
C	1003	GLU	-	expression tag	UNP Q66929
C	1004	GLU	-	expression tag	UNP Q66929
C	1005	LEU	-	expression tag	UNP Q66929
C	1006	ARG	-	expression tag	UNP Q66929
C	1007	ARG	-	expression tag	UNP Q66929
C	1008	ARG	-	expression tag	UNP Q66929
C	1009	LEU	-	expression tag	UNP Q66929
C	1010	THR	-	expression tag	UNP Q66929
C	1011	GLU	-	expression tag	UNP Q66929
D	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
D	999	SER	-	expression tag	UNP Q66929
D	1000	ARG	-	expression tag	UNP Q66929
D	1001	LEU	-	expression tag	UNP Q66929
D	1002	GLU	-	expression tag	UNP Q66929
D	1003	GLU	-	expression tag	UNP Q66929
D	1004	GLU	-	expression tag	UNP Q66929
D	1005	LEU	-	expression tag	UNP Q66929
D	1006	ARG	-	expression tag	UNP Q66929
D	1007	ARG	-	expression tag	UNP Q66929
D	1008	ARG	-	expression tag	UNP Q66929
D	1009	LEU	-	expression tag	UNP Q66929
D	1010	THR	-	expression tag	UNP Q66929
D	1011	GLU	-	expression tag	UNP Q66929
E	897	LEU	MET	conflict	UNP Q66929
E	999	SER	-	expression tag	UNP Q66929
E	1000	ARG	-	expression tag	UNP Q66929
E	1001	LEU	-	expression tag	UNP Q66929
E	1002	GLU	-	expression tag	UNP Q66929
E	1003	GLU	-	expression tag	UNP Q66929
E	1004	GLU	-	expression tag	UNP Q66929
E	1005	LEU	-	expression tag	UNP Q66929
E	1006	ARG	-	expression tag	UNP Q66929
E	1007	ARG	-	expression tag	UNP Q66929
E	1008	ARG	-	expression tag	UNP Q66929
E	1009	LEU	-	expression tag	UNP Q66929
E	1010	THR	-	expression tag	UNP Q66929
E	1011	GLU	-	expression tag	UNP Q66929
F	897	LEU	MET	conflict	UNP Q66929
F	999	SER	-	expression tag	UNP Q66929
F	1000	ARG	-	expression tag	UNP Q66929
F	1001	LEU	-	expression tag	UNP Q66929
F	1002	GLU	-	expression tag	UNP Q66929
F	1003	GLU	-	expression tag	UNP Q66929
F	1004	GLU	-	expression tag	UNP Q66929
F	1005	LEU	-	expression tag	UNP Q66929
F	1006	ARG	-	expression tag	UNP Q66929
F	1007	ARG	-	expression tag	UNP Q66929
F	1008	ARG	-	expression tag	UNP Q66929
F	1009	LEU	-	expression tag	UNP Q66929
F	1010	THR	-	expression tag	UNP Q66929
F	1011	GLU	-	expression tag	UNP Q66929
G	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
G	999	SER	-	expression tag	UNP Q66929
G	1000	ARG	-	expression tag	UNP Q66929
G	1001	LEU	-	expression tag	UNP Q66929
G	1002	GLU	-	expression tag	UNP Q66929
G	1003	GLU	-	expression tag	UNP Q66929
G	1004	GLU	-	expression tag	UNP Q66929
G	1005	LEU	-	expression tag	UNP Q66929
G	1006	ARG	-	expression tag	UNP Q66929
G	1007	ARG	-	expression tag	UNP Q66929
G	1008	ARG	-	expression tag	UNP Q66929
G	1009	LEU	-	expression tag	UNP Q66929
G	1010	THR	-	expression tag	UNP Q66929
G	1011	GLU	-	expression tag	UNP Q66929
H	897	LEU	MET	conflict	UNP Q66929
H	999	SER	-	expression tag	UNP Q66929
H	1000	ARG	-	expression tag	UNP Q66929
H	1001	LEU	-	expression tag	UNP Q66929
H	1002	GLU	-	expression tag	UNP Q66929
H	1003	GLU	-	expression tag	UNP Q66929
H	1004	GLU	-	expression tag	UNP Q66929
H	1005	LEU	-	expression tag	UNP Q66929
H	1006	ARG	-	expression tag	UNP Q66929
H	1007	ARG	-	expression tag	UNP Q66929
H	1008	ARG	-	expression tag	UNP Q66929
H	1009	LEU	-	expression tag	UNP Q66929
H	1010	THR	-	expression tag	UNP Q66929
H	1011	GLU	-	expression tag	UNP Q66929
I	897	LEU	MET	conflict	UNP Q66929
I	999	SER	-	expression tag	UNP Q66929
I	1000	ARG	-	expression tag	UNP Q66929
I	1001	LEU	-	expression tag	UNP Q66929
I	1002	GLU	-	expression tag	UNP Q66929
I	1003	GLU	-	expression tag	UNP Q66929
I	1004	GLU	-	expression tag	UNP Q66929
I	1005	LEU	-	expression tag	UNP Q66929
I	1006	ARG	-	expression tag	UNP Q66929
I	1007	ARG	-	expression tag	UNP Q66929
I	1008	ARG	-	expression tag	UNP Q66929
I	1009	LEU	-	expression tag	UNP Q66929
I	1010	THR	-	expression tag	UNP Q66929
I	1011	GLU	-	expression tag	UNP Q66929
J	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
J	999	SER	-	expression tag	UNP Q66929
J	1000	ARG	-	expression tag	UNP Q66929
J	1001	LEU	-	expression tag	UNP Q66929
J	1002	GLU	-	expression tag	UNP Q66929
J	1003	GLU	-	expression tag	UNP Q66929
J	1004	GLU	-	expression tag	UNP Q66929
J	1005	LEU	-	expression tag	UNP Q66929
J	1006	ARG	-	expression tag	UNP Q66929
J	1007	ARG	-	expression tag	UNP Q66929
J	1008	ARG	-	expression tag	UNP Q66929
J	1009	LEU	-	expression tag	UNP Q66929
J	1010	THR	-	expression tag	UNP Q66929
J	1011	GLU	-	expression tag	UNP Q66929
K	897	LEU	MET	conflict	UNP Q66929
K	999	SER	-	expression tag	UNP Q66929
K	1000	ARG	-	expression tag	UNP Q66929
K	1001	LEU	-	expression tag	UNP Q66929
K	1002	GLU	-	expression tag	UNP Q66929
K	1003	GLU	-	expression tag	UNP Q66929
K	1004	GLU	-	expression tag	UNP Q66929
K	1005	LEU	-	expression tag	UNP Q66929
K	1006	ARG	-	expression tag	UNP Q66929
K	1007	ARG	-	expression tag	UNP Q66929
K	1008	ARG	-	expression tag	UNP Q66929
K	1009	LEU	-	expression tag	UNP Q66929
K	1010	THR	-	expression tag	UNP Q66929
K	1011	GLU	-	expression tag	UNP Q66929
A	897	LEU	MET	conflict	UNP Q66929
A	999	SER	-	expression tag	UNP Q66929
A	1000	ARG	-	expression tag	UNP Q66929
A	1001	LEU	-	expression tag	UNP Q66929
A	1002	GLU	-	expression tag	UNP Q66929
A	1003	GLU	-	expression tag	UNP Q66929
A	1004	GLU	-	expression tag	UNP Q66929
A	1005	LEU	-	expression tag	UNP Q66929
A	1006	ARG	-	expression tag	UNP Q66929
A	1007	ARG	-	expression tag	UNP Q66929
A	1008	ARG	-	expression tag	UNP Q66929
A	1009	LEU	-	expression tag	UNP Q66929
A	1010	THR	-	expression tag	UNP Q66929
A	1011	GLU	-	expression tag	UNP Q66929
L	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
L	999	SER	-	expression tag	UNP Q66929
L	1000	ARG	-	expression tag	UNP Q66929
L	1001	LEU	-	expression tag	UNP Q66929
L	1002	GLU	-	expression tag	UNP Q66929
L	1003	GLU	-	expression tag	UNP Q66929
L	1004	GLU	-	expression tag	UNP Q66929
L	1005	LEU	-	expression tag	UNP Q66929
L	1006	ARG	-	expression tag	UNP Q66929
L	1007	ARG	-	expression tag	UNP Q66929
L	1008	ARG	-	expression tag	UNP Q66929
L	1009	LEU	-	expression tag	UNP Q66929
L	1010	THR	-	expression tag	UNP Q66929
L	1011	GLU	-	expression tag	UNP Q66929
M	897	LEU	MET	conflict	UNP Q66929
M	999	SER	-	expression tag	UNP Q66929
M	1000	ARG	-	expression tag	UNP Q66929
M	1001	LEU	-	expression tag	UNP Q66929
M	1002	GLU	-	expression tag	UNP Q66929
M	1003	GLU	-	expression tag	UNP Q66929
M	1004	GLU	-	expression tag	UNP Q66929
M	1005	LEU	-	expression tag	UNP Q66929
M	1006	ARG	-	expression tag	UNP Q66929
M	1007	ARG	-	expression tag	UNP Q66929
M	1008	ARG	-	expression tag	UNP Q66929
M	1009	LEU	-	expression tag	UNP Q66929
M	1010	THR	-	expression tag	UNP Q66929
M	1011	GLU	-	expression tag	UNP Q66929
N	897	LEU	MET	conflict	UNP Q66929
N	999	SER	-	expression tag	UNP Q66929
N	1000	ARG	-	expression tag	UNP Q66929
N	1001	LEU	-	expression tag	UNP Q66929
N	1002	GLU	-	expression tag	UNP Q66929
N	1003	GLU	-	expression tag	UNP Q66929
N	1004	GLU	-	expression tag	UNP Q66929
N	1005	LEU	-	expression tag	UNP Q66929
N	1006	ARG	-	expression tag	UNP Q66929
N	1007	ARG	-	expression tag	UNP Q66929
N	1008	ARG	-	expression tag	UNP Q66929
N	1009	LEU	-	expression tag	UNP Q66929
N	1010	THR	-	expression tag	UNP Q66929
N	1011	GLU	-	expression tag	UNP Q66929
O	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
O	999	SER	-	expression tag	UNP Q66929
O	1000	ARG	-	expression tag	UNP Q66929
O	1001	LEU	-	expression tag	UNP Q66929
O	1002	GLU	-	expression tag	UNP Q66929
O	1003	GLU	-	expression tag	UNP Q66929
O	1004	GLU	-	expression tag	UNP Q66929
O	1005	LEU	-	expression tag	UNP Q66929
O	1006	ARG	-	expression tag	UNP Q66929
O	1007	ARG	-	expression tag	UNP Q66929
O	1008	ARG	-	expression tag	UNP Q66929
O	1009	LEU	-	expression tag	UNP Q66929
O	1010	THR	-	expression tag	UNP Q66929
O	1011	GLU	-	expression tag	UNP Q66929
P	897	LEU	MET	conflict	UNP Q66929
P	999	SER	-	expression tag	UNP Q66929
P	1000	ARG	-	expression tag	UNP Q66929
P	1001	LEU	-	expression tag	UNP Q66929
P	1002	GLU	-	expression tag	UNP Q66929
P	1003	GLU	-	expression tag	UNP Q66929
P	1004	GLU	-	expression tag	UNP Q66929
P	1005	LEU	-	expression tag	UNP Q66929
P	1006	ARG	-	expression tag	UNP Q66929
P	1007	ARG	-	expression tag	UNP Q66929
P	1008	ARG	-	expression tag	UNP Q66929
P	1009	LEU	-	expression tag	UNP Q66929
P	1010	THR	-	expression tag	UNP Q66929
P	1011	GLU	-	expression tag	UNP Q66929
Q	897	LEU	MET	conflict	UNP Q66929
Q	999	SER	-	expression tag	UNP Q66929
Q	1000	ARG	-	expression tag	UNP Q66929
Q	1001	LEU	-	expression tag	UNP Q66929
Q	1002	GLU	-	expression tag	UNP Q66929
Q	1003	GLU	-	expression tag	UNP Q66929
Q	1004	GLU	-	expression tag	UNP Q66929
Q	1005	LEU	-	expression tag	UNP Q66929
Q	1006	ARG	-	expression tag	UNP Q66929
Q	1007	ARG	-	expression tag	UNP Q66929
Q	1008	ARG	-	expression tag	UNP Q66929
Q	1009	LEU	-	expression tag	UNP Q66929
Q	1010	THR	-	expression tag	UNP Q66929
Q	1011	GLU	-	expression tag	UNP Q66929
R	897	LEU	MET	conflict	UNP Q66929

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Chain	Residue	Modelled	Actual	Comment	Reference
R	999	SER	-	expression tag	UNP Q66929
R	1000	ARG	-	expression tag	UNP Q66929
R	1001	LEU	-	expression tag	UNP Q66929
R	1002	GLU	-	expression tag	UNP Q66929
R	1003	GLU	-	expression tag	UNP Q66929
R	1004	GLU	-	expression tag	UNP Q66929
R	1005	LEU	-	expression tag	UNP Q66929
R	1006	ARG	-	expression tag	UNP Q66929
R	1007	ARG	-	expression tag	UNP Q66929
R	1008	ARG	-	expression tag	UNP Q66929
R	1009	LEU	-	expression tag	UNP Q66929
R	1010	THR	-	expression tag	UNP Q66929
R	1011	GLU	-	expression tag	UNP Q66929
S	897	LEU	MET	conflict	UNP Q66929
S	999	SER	-	expression tag	UNP Q66929
S	1000	ARG	-	expression tag	UNP Q66929
S	1001	LEU	-	expression tag	UNP Q66929
S	1002	GLU	-	expression tag	UNP Q66929
S	1003	GLU	-	expression tag	UNP Q66929
S	1004	GLU	-	expression tag	UNP Q66929
S	1005	LEU	-	expression tag	UNP Q66929
S	1006	ARG	-	expression tag	UNP Q66929
S	1007	ARG	-	expression tag	UNP Q66929
S	1008	ARG	-	expression tag	UNP Q66929
S	1009	LEU	-	expression tag	UNP Q66929
S	1010	THR	-	expression tag	UNP Q66929
S	1011	GLU	-	expression tag	UNP Q66929
T	897	LEU	MET	conflict	UNP Q66929
T	999	SER	-	expression tag	UNP Q66929
T	1000	ARG	-	expression tag	UNP Q66929
T	1001	LEU	-	expression tag	UNP Q66929
T	1002	GLU	-	expression tag	UNP Q66929
T	1003	GLU	-	expression tag	UNP Q66929
T	1004	GLU	-	expression tag	UNP Q66929
T	1005	LEU	-	expression tag	UNP Q66929
T	1006	ARG	-	expression tag	UNP Q66929
T	1007	ARG	-	expression tag	UNP Q66929
T	1008	ARG	-	expression tag	UNP Q66929
T	1009	LEU	-	expression tag	UNP Q66929
T	1010	THR	-	expression tag	UNP Q66929
T	1011	GLU	-	expression tag	UNP Q66929
U	897	LEU	MET	conflict	UNP Q66929

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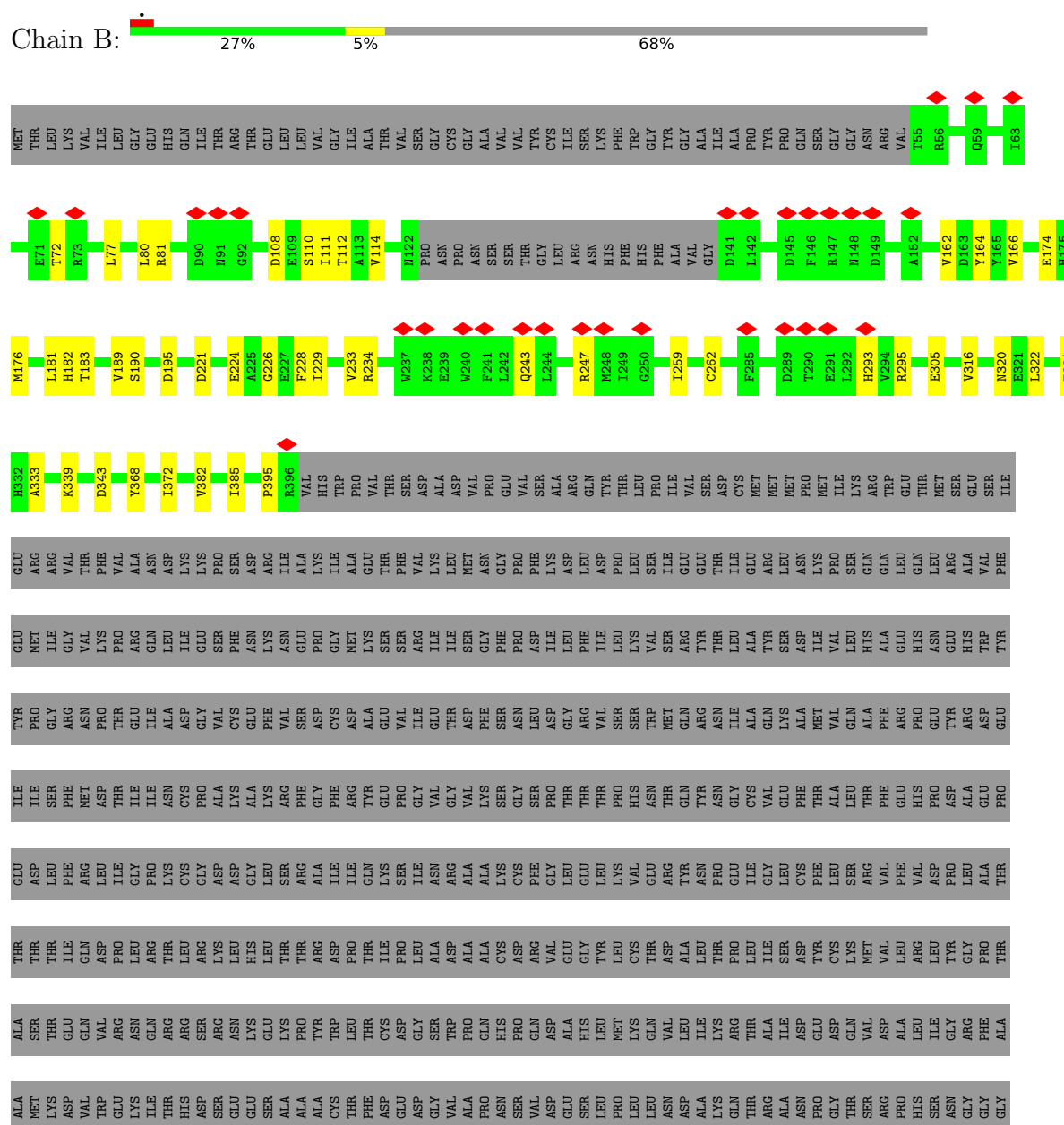
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Chain	Residue	Modelled	Actual	Comment	Reference
U	999	SER	-	expression tag	UNP Q66929
U	1000	ARG	-	expression tag	UNP Q66929
U	1001	LEU	-	expression tag	UNP Q66929
U	1002	GLU	-	expression tag	UNP Q66929
U	1003	GLU	-	expression tag	UNP Q66929
U	1004	GLU	-	expression tag	UNP Q66929
U	1005	LEU	-	expression tag	UNP Q66929
U	1006	ARG	-	expression tag	UNP Q66929
U	1007	ARG	-	expression tag	UNP Q66929
U	1008	ARG	-	expression tag	UNP Q66929
U	1009	LEU	-	expression tag	UNP Q66929
U	1010	THR	-	expression tag	UNP Q66929
U	1011	GLU	-	expression tag	UNP Q66929
V	897	LEU	MET	conflict	UNP Q66929
V	999	SER	-	expression tag	UNP Q66929
V	1000	ARG	-	expression tag	UNP Q66929
V	1001	LEU	-	expression tag	UNP Q66929
V	1002	GLU	-	expression tag	UNP Q66929
V	1003	GLU	-	expression tag	UNP Q66929
V	1004	GLU	-	expression tag	UNP Q66929
V	1005	LEU	-	expression tag	UNP Q66929
V	1006	ARG	-	expression tag	UNP Q66929
V	1007	ARG	-	expression tag	UNP Q66929
V	1008	ARG	-	expression tag	UNP Q66929
V	1009	LEU	-	expression tag	UNP Q66929
V	1010	THR	-	expression tag	UNP Q66929
V	1011	GLU	-	expression tag	UNP Q66929

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

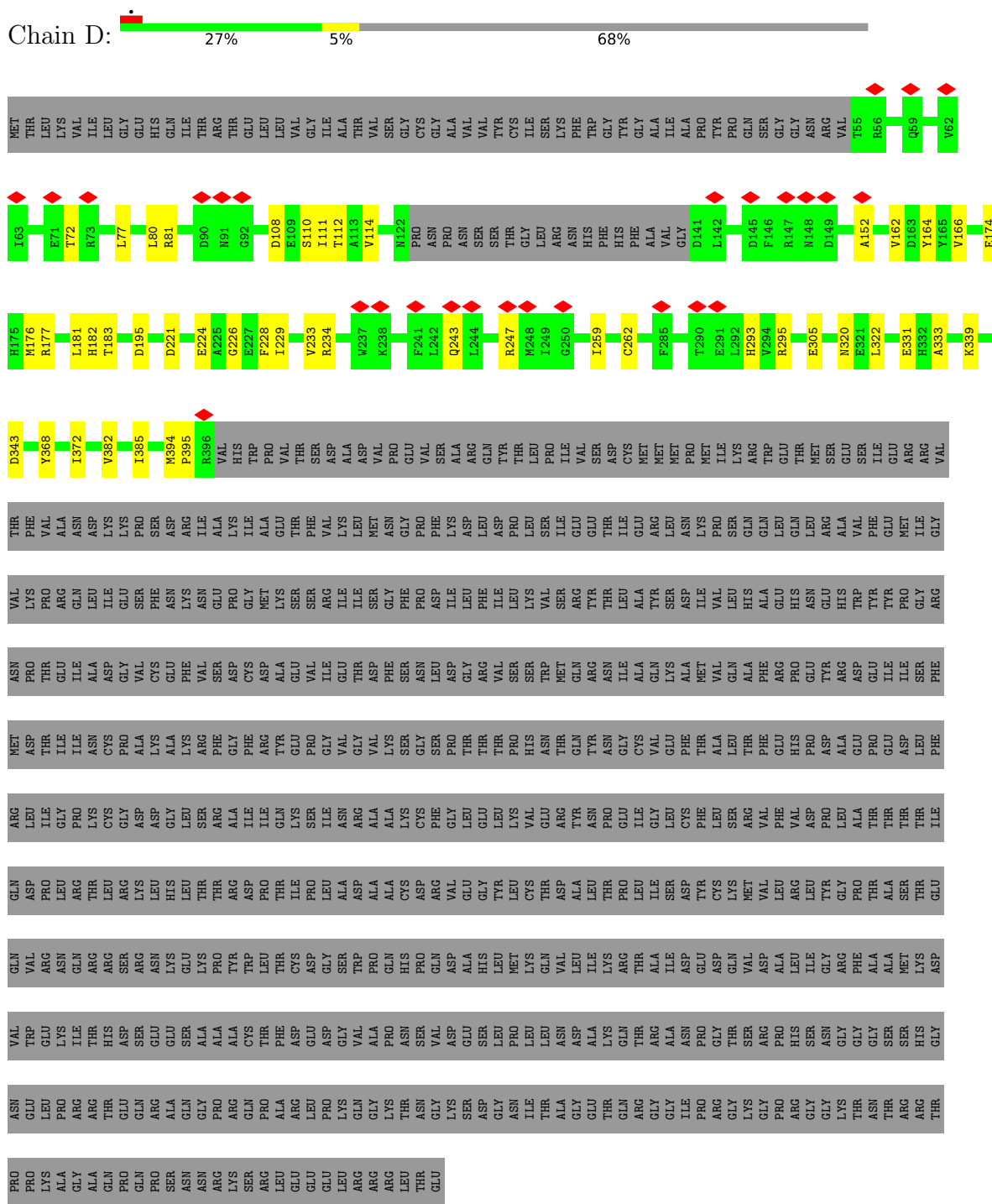
• Molecule 1: RNA-directed RNA polymerase



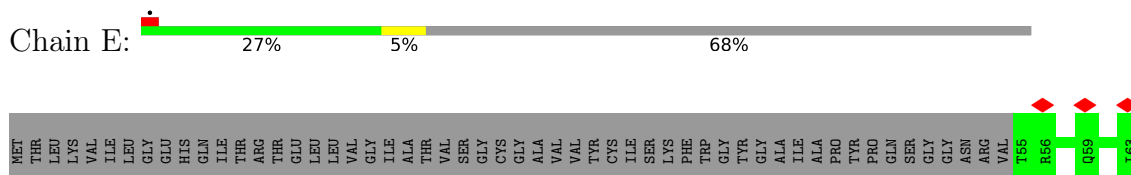
- Molecule 1: RNA-directed RNA polymerase

[illegible]

- Molecule 1: RNA-directed RNA polymerase



- Molecule 1: RNA-directed RNA polymerase



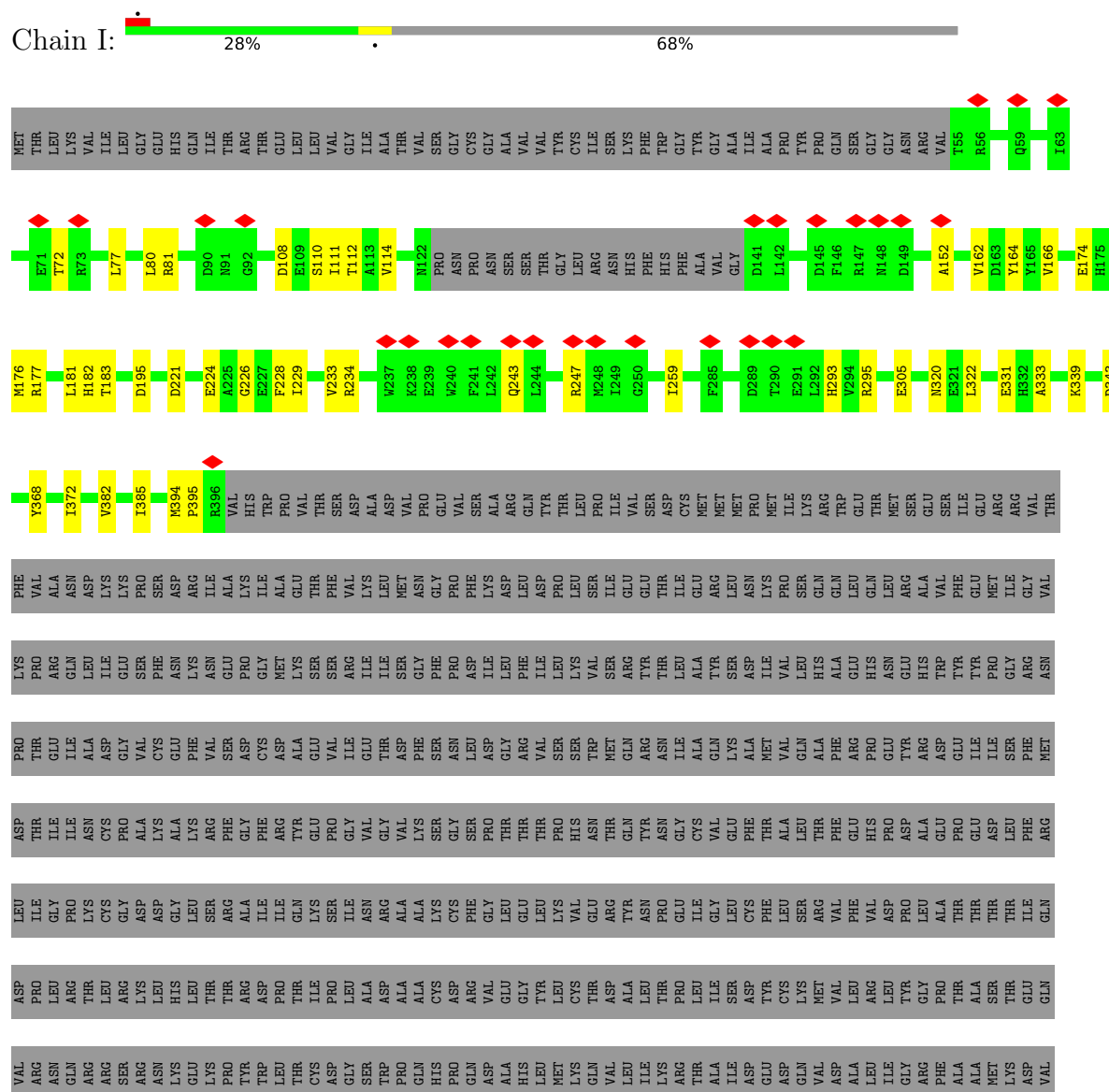
LYS	LEU	GLU	ARG	PRO	ILE	THR	THR	PRO
ALA	PRO	LYS	ASN	LEU	GLY	ILE	GLU	ARG
GLY	ARG	ILE	GLN	ARG	PRO	ILE	ILE	GLN
ALA	THR	HIS	ARG	THR	LYS	ASN	ALA	LEU
GLN	THR	HIS	ARG	LEU	CYS	ASN	GLY	ILE
PRO	GLU	ASP	SER	ARG	GLY	PRO	GLY	GLU
GLN	GLN	SER	ARG	LYS	ASP	ALA	VAL	SER
PRO	ARG	GLU	ASN	LEU	ASN	LYS	CYS	PHE
SER	ALA	GLU	LYS	HIS	GLY	ALA	GLU	ASN
ASN	GLN	SER	GLU	LEU	LEU	LYS	PHE	LYS
ASN	GLY	ALA	LYS	THR	SER	ARG	VAL	ASN
ARG	PRO	ALA	PRO	THR	ARG	PHE	SER	GLU
LYS	ARG	ALA	TYP	ARG	GLY	GLY	ASP	PRO
SER	GLN	CYS	TRP	ASP	ILE	PHE	CYS	GLY
ARG	PRO	THR	LEU	PRO	ILE	ARG	ASP	MET
LEU	ALA	PHE	THR	THR	GLN	TYP	ALA	LYS
GLU	ARG	ASP	CYS	ILE	LYS	GLY	GLU	SER
GLU	LEU	GLU	ASP	PRO	SER	PRO	VAL	SER
LEU	PRO	ASP	GLY	LEU	ILE	GLY	ILE	ARG
LEU	LYS	GLY	SER	ALA	ASN	VAL	GLU	ILE
ARG	GLN	VAL	TRP	ASP	ARG	GLY	THR	ILE
ARG	GLY	ALA	PRO	ALA	ALA	VAL	ASP	SER
ARG	LYS	PRO	GLN	ALA	LYS	SER	PHE	GLY
LEU	THR	ASN	HIS	CYS	LYS	SER	SER	PHE
THR	ASN	SER	PRO	ASP	CYS	GLY	ASN	PHE
THR	VAL	VAL	GLN	ARG	PHE	SER	LEU	ARG
GLY	GLY	LYS	ASP	VAL	GLY	PRO	ASP	ILE
SER	SER	GLU	ALA	GLU	LEU	THR	GLY	PHE
ASP	ASP	SER	HIS	GLY	GLU	THR	ARG	PHE
GLY	GLY	LEU	LEU	TYP	LEU	THR	VAL	ILE
ASN	ASN	PRO	MET	LEU	LYS	PRO	SER	ILE
ILE	ILE	LEU	LYS	CYS	VAL	HIS	LYS	LEU
THR	THR	LEU	GLN	THR	GLU	ASN	TRP	VAL
ALA	ALA	ASN	VAL	ASP	ARG	THR	MET	SER
GLY	GLY	ASP	LEU	ALA	TYP	GLN	GLN	ARG
ILE	ILE	ASN	ASP	SER	LEU	PHE	ALA	ASP
PRO	PRO	PRO	GLU	TYP	PHE	THR	MET	ILE
ARG	ARG	GLY	ASP	CYS	LEU	ALA	VAL	VAL
GLY	GLY	THR	GLN	LYS	SER	LEU	GLN	LEU
LYS	LYS	ARG	VAL	MET	ARG	THR	ALA	ALA
PRO	PRO	PRO	ALA	VAL	VAL	PHE	PHE	ALA
ARG	ARG	HIS	LEU	ARG	VAL	GLU	ARG	GLU
GLY	GLY	SER	ILE	THR	VAL	HIS	PRO	HIS
LYS	LYS	ASN	GLY	TYP	PRO	ASP	GLU	ASN
THR	THR	GLY	ARG	TYP	ALA	ALA	TYP	GLN
ASN	ASN	GLY	THR	PRO	THR	GLU	ASP	THR
PRO	PRO	GLY	GLY	GLY	LEU	THR	GLU	ARG
THR	THR	GLY	PHE	PRO	ALA	PRO	GLU	ASN
ASN	THR	GLY	ALA	THR	THR	PRO	GLU	PRO
PRO	THR	SER	MET	ALA	THR	ASP	ILE	THR
ARG	ARG	HIS	LYS	THR	THR	PHE	SER	ARG
THR	THR	GLY	ASP	GLU	ILE	ARG	PHE	THR
PRO	PRO	ASN	VAL	GLN	GLN	LEU	MET	PRO
THR	THR	GLU	THR	VAL	ASP	ARG	THR	ASN

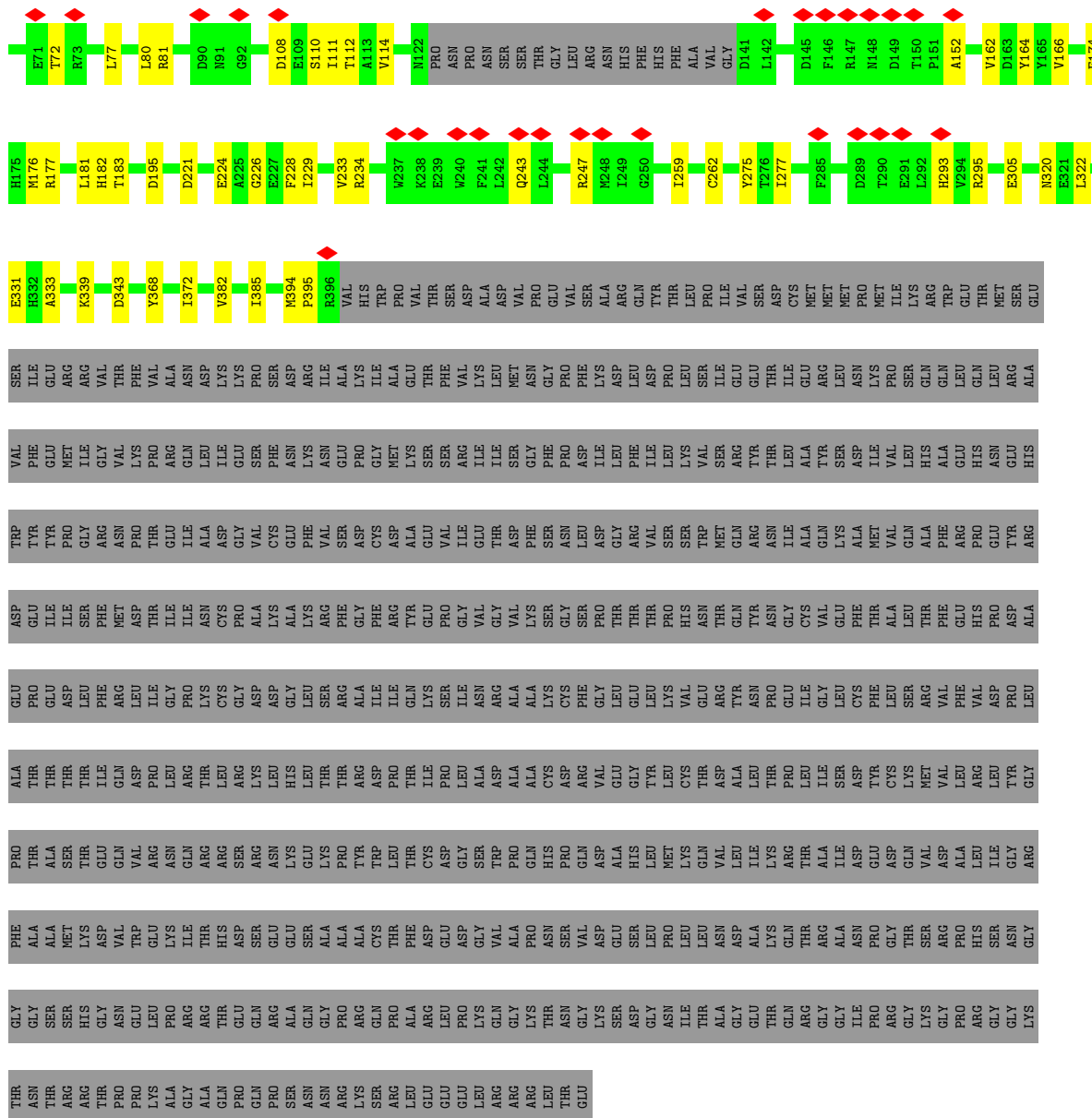
- Molecule 1: RNA-directed RNA polymerase

[illegible]

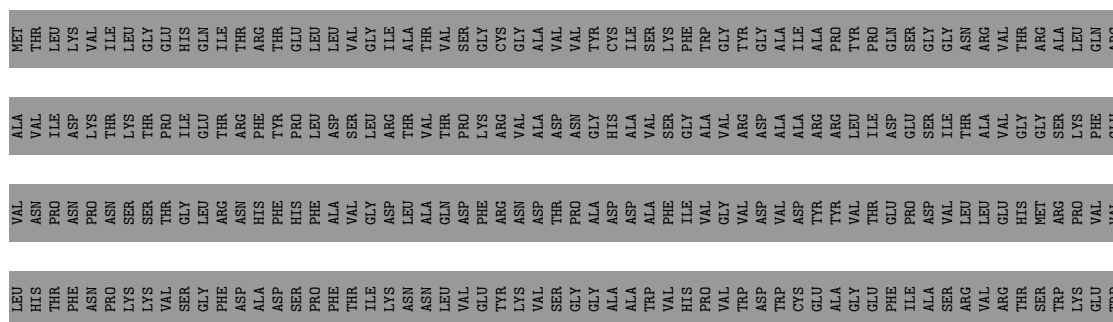
[illegible]

- Molecule 1: RNA-directed RNA polymerase





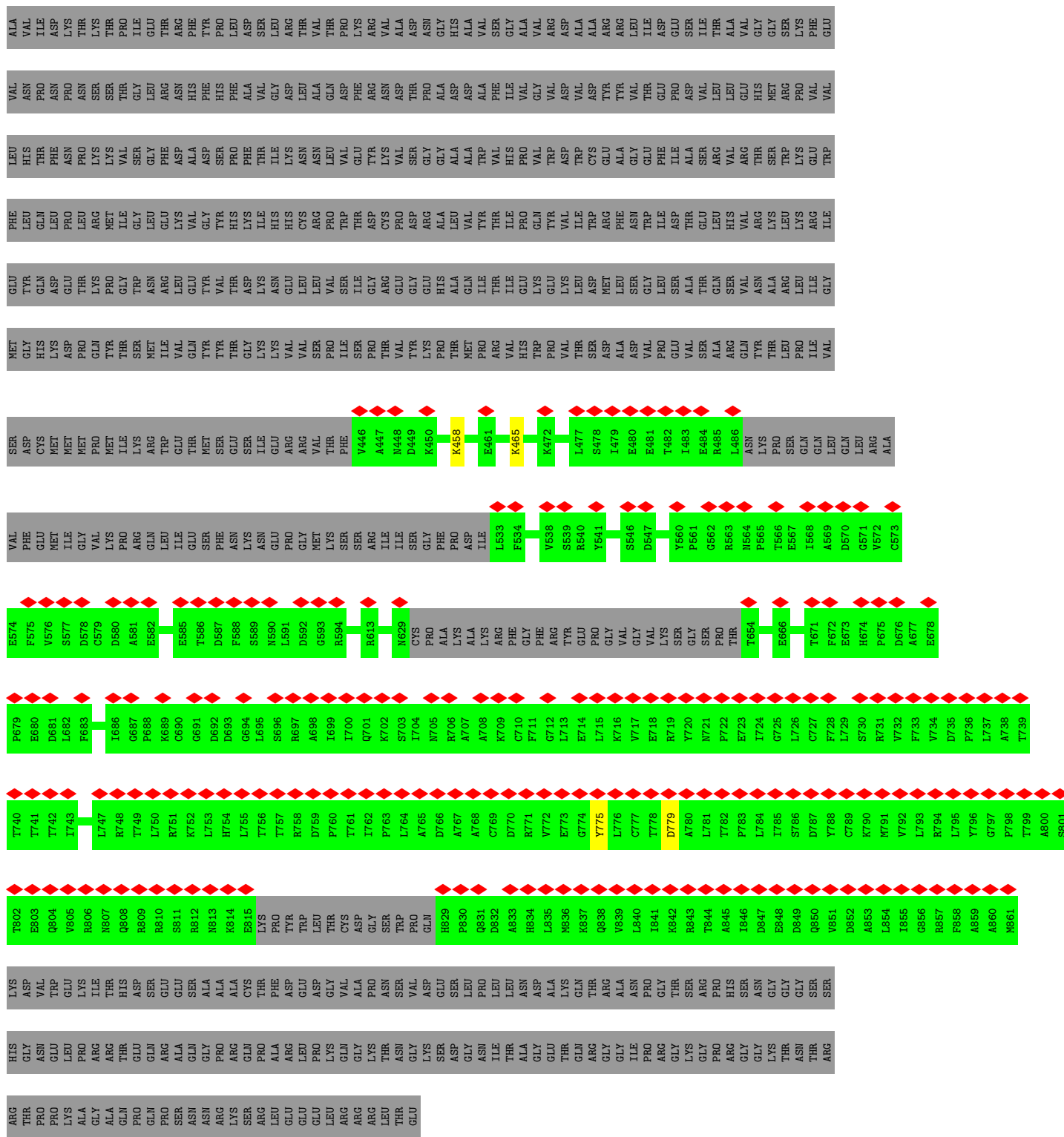
- Molecule 1: RNA-directed RNA polymerase



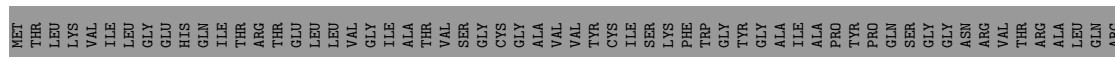
ASN	THR	ARG	ARG	THR	PRO	PRO	LYS	ALA	GLY	ALA	GLN	PRO	GLN	PRO	PRO	SER	SER	ASN	ASN	ASN	ARG	LYS	SER	SER	ARG	LEU	GLU	GLU	GLU	LEU	ARG	ARG	ARG	LEU	THR	GLU
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

Chain P: 21% 33% 67%

MET	THR	LEU	LEU	LYS	VAL	ILE	LEU	GLY	GLU	GLN	HIS	THR	THR	ARG	THR	GLU	LEU	LEU	VAL	GLY	ILE	ALA	ALA	THR	THR	VAL	VAL	VAL	TYR	CYS	CYS	ILE	ILE	LYS	PHE	TRP	GLY	GLY	TYR	TYR	PRO	PRO	GLN	GLN	SER	ALA	ALA	ILE	ALA	ALA	PRO	PRO	TYR	TYR	ASN	ASN	ARG	VAL	THR	THR	ARG	ALA	ALA	LEU	LEU	GLN	PC
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	----



- Molecule 1: RNA-directed RNA polymerase

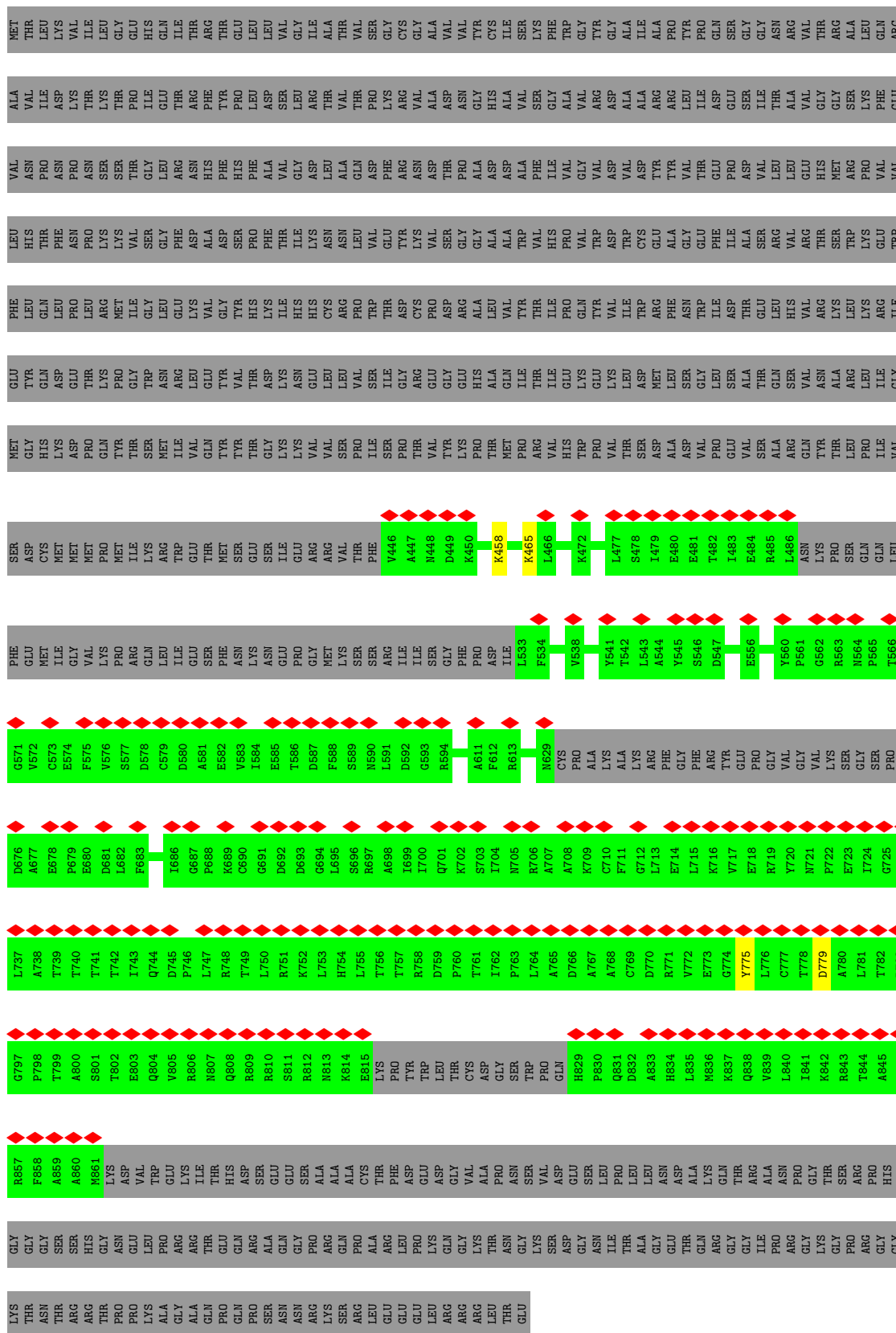




- Molecule 1: RNA-directed RNA polymerase

[illegible]





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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C11	Depositor
Number of particles used	13701	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	100	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	32.812	Depositor
Minimum map value	-20.300	Depositor
Average map value	-0.008	Depositor
Map value standard deviation	0.592	Depositor
Recommended contour level	3.09	Depositor
Map size (Å)	479.808, 479.808, 479.808	wwPDB
Map dimensions	576, 576, 576	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.833, 0.833, 0.833	Depositor

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.13	0/2688	0.31	0/3655
1	B	0.13	0/2688	0.31	0/3655
1	C	0.14	0/2688	0.31	0/3655
1	D	0.13	0/2688	0.31	0/3655
1	E	0.13	0/2688	0.31	0/3655
1	F	0.13	0/2688	0.31	0/3655
1	G	0.14	0/2688	0.31	0/3655
1	H	0.13	0/2688	0.31	0/3655
1	I	0.13	0/2688	0.31	0/3655
1	J	0.13	0/2688	0.31	0/3655
1	K	0.13	0/2688	0.31	0/3655
1	L	0.08	0/1409	0.27	0/1766
1	M	0.08	0/1409	0.27	0/1766
1	N	0.08	0/1409	0.27	0/1766
1	O	0.08	0/1409	0.27	0/1766
1	P	0.08	0/1409	0.27	0/1766
1	Q	0.08	0/1409	0.27	0/1766
1	R	0.08	0/1409	0.27	0/1766
1	S	0.08	0/1409	0.27	0/1766
1	T	0.08	0/1409	0.27	0/1766
1	U	0.08	0/1409	0.27	0/1766
1	V	0.08	0/1409	0.27	0/1766
All	All	0.12	0/45067	0.30	0/59631

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2620	2603	2600	58	0
1	B	2620	2603	2600	54	0
1	C	2620	2603	2600	54	0
1	D	2620	2603	2600	55	0
1	E	2620	2603	2600	55	0
1	F	2620	2603	2600	59	0
1	G	2620	2603	2600	55	0
1	H	2620	2603	2600	55	0
1	I	2620	2603	2600	53	0
1	J	2620	2603	2600	56	0
1	K	2620	2603	2600	55	0
1	L	1411	0	464	4	0
1	M	1411	0	464	3	0
1	N	1411	0	464	3	0
1	O	1411	0	464	3	0
1	P	1411	0	464	4	0
1	Q	1411	0	464	3	0
1	R	1411	0	464	3	0
1	S	1411	0	464	3	0
1	T	1411	0	464	3	0
1	U	1411	0	464	4	0
1	V	1411	0	464	3	0
All	All	44341	28633	33704	603	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (603) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:GLU:OE2	1:B:234:ARG:HA	1.20	1.35
1:J:174:GLU:OE2	1:J:234:ARG:HA	1.20	1.35
1:C:174:GLU:OE2	1:C:234:ARG:HA	1.20	1.34
1:I:174:GLU:OE2	1:I:234:ARG:HA	1.20	1.34
1:G:174:GLU:OE2	1:G:234:ARG:HA	1.20	1.34

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:174:GLU:OE2	1:F:234:ARG:HA	1.20	1.33
1:E:174:GLU:OE2	1:E:234:ARG:HA	1.20	1.30
1:K:174:GLU:OE2	1:K:234:ARG:HA	1.20	1.29
1:A:174:GLU:OE2	1:A:234:ARG:HA	1.20	1.28
1:D:174:GLU:OE2	1:D:234:ARG:HA	1.20	1.28
1:H:174:GLU:OE2	1:H:234:ARG:HA	1.20	1.27
1:K:77:LEU:HD23	1:K:80:LEU:CD1	1.77	1.15
1:A:77:LEU:HD23	1:A:80:LEU:CD1	1.77	1.15
1:D:77:LEU:HD23	1:D:80:LEU:CD1	1.77	1.14
1:G:77:LEU:HD23	1:G:80:LEU:CD1	1.77	1.14
1:E:77:LEU:HD23	1:E:80:LEU:CD1	1.77	1.14
1:J:77:LEU:HD23	1:J:80:LEU:CD1	1.77	1.14
1:H:77:LEU:HD23	1:H:80:LEU:CD1	1.77	1.14
1:F:77:LEU:HD23	1:F:80:LEU:CD1	1.77	1.13
1:C:77:LEU:HD23	1:C:80:LEU:HD12	1.13	1.13
1:C:77:LEU:HD23	1:C:80:LEU:CD1	1.77	1.13
1:H:77:LEU:HD23	1:H:80:LEU:HD12	1.13	1.13
1:G:77:LEU:HD23	1:G:80:LEU:HD12	1.13	1.13
1:B:77:LEU:HD23	1:B:80:LEU:CD1	1.77	1.13
1:I:77:LEU:HD23	1:I:80:LEU:CD1	1.77	1.12
1:D:77:LEU:HD23	1:D:80:LEU:HD12	1.13	1.12
1:B:77:LEU:HD23	1:B:80:LEU:HD12	1.13	1.11
1:I:77:LEU:HD23	1:I:80:LEU:HD12	1.13	1.11
1:A:77:LEU:HD23	1:A:80:LEU:HD12	1.13	1.10
1:F:77:LEU:HD23	1:F:80:LEU:HD12	1.13	1.10
1:E:77:LEU:HD23	1:E:80:LEU:HD12	1.13	1.09
1:J:77:LEU:HD23	1:J:80:LEU:HD12	1.13	1.08
1:K:77:LEU:HD23	1:K:80:LEU:HD12	1.13	1.07
1:K:77:LEU:CD2	1:K:80:LEU:CD1	2.37	1.02
1:E:77:LEU:CD2	1:E:80:LEU:CD1	2.37	1.02
1:F:77:LEU:CD2	1:F:80:LEU:CD1	2.37	1.02
1:J:77:LEU:CD2	1:J:80:LEU:CD1	2.37	1.02
1:A:77:LEU:CD2	1:A:80:LEU:CD1	2.37	1.02
1:C:77:LEU:CD2	1:C:80:LEU:CD1	2.37	1.02
1:H:77:LEU:CD2	1:H:80:LEU:CD1	2.37	1.02
1:D:77:LEU:CD2	1:D:80:LEU:CD1	2.37	1.01
1:G:77:LEU:CD2	1:G:80:LEU:CD1	2.37	1.01
1:I:77:LEU:CD2	1:I:80:LEU:CD1	2.37	1.01
1:B:77:LEU:CD2	1:B:80:LEU:CD1	2.37	1.00
1:B:174:GLU:OE2	1:B:234:ARG:CA	2.10	1.00
1:C:174:GLU:OE2	1:C:234:ARG:CA	2.10	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:174:GLU:OE2	1:D:234:ARG:CA	2.10	1.00
1:I:174:GLU:OE2	1:I:234:ARG:CA	2.10	1.00
1:H:174:GLU:OE2	1:H:234:ARG:CA	2.10	1.00
1:J:174:GLU:OE2	1:J:234:ARG:CA	2.10	1.00
1:G:174:GLU:OE2	1:G:234:ARG:CA	2.10	1.00
1:E:174:GLU:OE2	1:E:234:ARG:CA	2.10	0.99
1:A:174:GLU:OE2	1:A:234:ARG:CA	2.10	0.99
1:K:174:GLU:OE2	1:K:234:ARG:CA	2.10	0.99
1:F:174:GLU:OE2	1:F:234:ARG:CA	2.10	0.99
1:A:77:LEU:HD21	1:A:80:LEU:HG	1.54	0.89
1:C:77:LEU:HD21	1:C:80:LEU:HG	1.54	0.89
1:B:77:LEU:HD21	1:B:80:LEU:HG	1.55	0.89
1:H:77:LEU:HD21	1:H:80:LEU:HG	1.55	0.88
1:G:77:LEU:HD21	1:G:80:LEU:HG	1.54	0.88
1:I:77:LEU:HD21	1:I:80:LEU:HG	1.54	0.88
1:K:77:LEU:HD21	1:K:80:LEU:HG	1.55	0.88
1:J:77:LEU:HD21	1:J:80:LEU:HG	1.54	0.87
1:D:77:LEU:HD21	1:D:80:LEU:HG	1.55	0.87
1:F:77:LEU:HD21	1:F:80:LEU:HG	1.55	0.87
1:E:77:LEU:HD21	1:E:80:LEU:HG	1.55	0.85
1:G:77:LEU:CD2	1:G:80:LEU:HG	2.11	0.81
1:D:77:LEU:CD2	1:D:80:LEU:HG	2.11	0.81
1:I:77:LEU:CD2	1:I:80:LEU:HG	2.11	0.81
1:K:77:LEU:CD2	1:K:80:LEU:HG	2.11	0.81
1:A:77:LEU:CD2	1:A:80:LEU:HG	2.11	0.81
1:E:77:LEU:CD2	1:E:80:LEU:HG	2.11	0.81
1:B:77:LEU:CD2	1:B:80:LEU:HG	2.11	0.81
1:F:77:LEU:CD2	1:F:80:LEU:HG	2.11	0.80
1:J:77:LEU:CD2	1:J:80:LEU:HG	2.11	0.80
1:C:77:LEU:CD2	1:C:80:LEU:HG	2.11	0.80
1:D:81:ARG:HD2	1:I:395:PRO:HG3	1.64	0.80
1:H:77:LEU:CD2	1:H:80:LEU:HG	2.11	0.80
1:H:174:GLU:HG3	1:H:233:VAL:C	2.08	0.79
1:C:81:ARG:HD2	1:H:395:PRO:HG3	1.64	0.79
1:C:174:GLU:HG3	1:C:233:VAL:C	2.08	0.79
1:G:174:GLU:HG3	1:G:233:VAL:C	2.08	0.79
1:B:174:GLU:HG3	1:B:233:VAL:C	2.08	0.79
1:A:174:GLU:HG3	1:A:233:VAL:C	2.08	0.78
1:B:81:ARG:HD2	1:G:395:PRO:HG3	1.65	0.78
1:J:174:GLU:HG3	1:J:233:VAL:C	2.08	0.78
1:D:174:GLU:HG3	1:D:233:VAL:C	2.08	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:174:GLU:HG3	1:F:233:VAL:C	2.08	0.78
1:K:174:GLU:HG3	1:K:233:VAL:C	2.08	0.78
1:E:174:GLU:HG3	1:E:233:VAL:C	2.08	0.78
1:I:174:GLU:HG3	1:I:233:VAL:C	2.08	0.78
1:F:395:PRO:HG3	1:A:81:ARG:HD2	1.68	0.76
1:C:77:LEU:CD2	1:C:80:LEU:HD12	2.01	0.76
1:G:77:LEU:HD21	1:G:80:LEU:CG	2.16	0.76
1:C:77:LEU:HD21	1:C:80:LEU:CG	2.16	0.75
1:H:77:LEU:HD21	1:H:80:LEU:CG	2.16	0.75
1:F:77:LEU:CD2	1:F:80:LEU:CG	2.65	0.75
1:K:77:LEU:HD21	1:K:80:LEU:CG	2.16	0.75
1:J:77:LEU:CD2	1:J:80:LEU:CG	2.65	0.75
1:B:77:LEU:HD21	1:B:80:LEU:CG	2.16	0.74
1:J:77:LEU:HD21	1:J:80:LEU:CG	2.16	0.74
1:A:77:LEU:HD21	1:A:80:LEU:CG	2.16	0.74
1:D:77:LEU:HD21	1:D:80:LEU:CG	2.16	0.74
1:K:77:LEU:CD2	1:K:80:LEU:CG	2.65	0.74
1:E:77:LEU:CD2	1:E:80:LEU:CG	2.65	0.74
1:B:77:LEU:CD2	1:B:80:LEU:CG	2.65	0.74
1:F:77:LEU:HD21	1:F:80:LEU:CG	2.16	0.74
1:H:77:LEU:CD2	1:H:80:LEU:CG	2.65	0.74
1:E:77:LEU:HD21	1:E:80:LEU:CG	2.16	0.74
1:I:77:LEU:CD2	1:I:80:LEU:CG	2.65	0.74
1:D:77:LEU:CD2	1:D:80:LEU:CG	2.65	0.74
1:I:77:LEU:HD21	1:I:80:LEU:CG	2.16	0.74
1:A:77:LEU:CD2	1:A:80:LEU:CG	2.65	0.74
1:C:77:LEU:CD2	1:C:80:LEU:CG	2.65	0.73
1:G:77:LEU:CD2	1:G:80:LEU:CG	2.65	0.73
1:D:395:PRO:HG3	1:J:81:ARG:HD2	1.69	0.73
1:G:77:LEU:HD11	1:G:110:SER:HB2	1.71	0.73
1:G:81:ARG:HD2	1:A:395:PRO:HG3	1.70	0.72
1:K:77:LEU:HD11	1:K:110:SER:HB2	1.71	0.72
1:E:77:LEU:HD11	1:E:110:SER:HB2	1.71	0.72
1:H:77:LEU:HD11	1:H:110:SER:HB2	1.71	0.72
1:J:77:LEU:HD11	1:J:110:SER:HB2	1.72	0.72
1:F:77:LEU:HD11	1:F:110:SER:HB2	1.71	0.72
1:A:77:LEU:HD11	1:A:110:SER:HB2	1.72	0.72
1:K:77:LEU:HD21	1:K:80:LEU:CD1	2.20	0.72
1:I:77:LEU:HD11	1:I:110:SER:HB2	1.71	0.72
1:B:77:LEU:HD11	1:B:110:SER:HB2	1.71	0.72
1:E:81:ARG:HD2	1:J:395:PRO:HG3	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:LEU:HD11	1:D:110:SER:HB2	1.71	0.72
1:F:77:LEU:HD21	1:F:80:LEU:CD1	2.20	0.72
1:G:77:LEU:CD2	1:G:80:LEU:HD12	2.01	0.71
1:C:77:LEU:HD11	1:C:110:SER:HB2	1.71	0.71
1:B:77:LEU:HD21	1:B:80:LEU:CD1	2.20	0.71
1:E:77:LEU:HD21	1:E:80:LEU:CD1	2.20	0.71
1:A:77:LEU:HD21	1:A:80:LEU:CD1	2.20	0.71
1:D:77:LEU:CD2	1:D:80:LEU:HD12	2.01	0.70
1:D:77:LEU:HD21	1:D:80:LEU:CD1	2.20	0.70
1:E:395:PRO:HG3	1:K:81:ARG:HD2	1.72	0.70
1:F:81:ARG:HD2	1:K:395:PRO:HG3	1.73	0.70
1:H:77:LEU:HD21	1:H:80:LEU:CD1	2.20	0.70
1:I:77:LEU:HD21	1:I:80:LEU:CD1	2.20	0.70
1:J:77:LEU:HD21	1:J:80:LEU:CD1	2.20	0.70
1:K:174:GLU:CG	1:K:234:ARG:N	2.55	0.70
1:F:174:GLU:CG	1:F:234:ARG:N	2.55	0.70
1:H:174:GLU:CG	1:H:234:ARG:N	2.55	0.70
1:J:174:GLU:CG	1:J:234:ARG:N	2.55	0.70
1:C:77:LEU:HD21	1:C:80:LEU:CD1	2.20	0.70
1:C:174:GLU:CG	1:C:234:ARG:N	2.55	0.70
1:G:174:GLU:CG	1:G:234:ARG:N	2.55	0.70
1:E:174:GLU:CG	1:E:234:ARG:N	2.55	0.69
1:H:77:LEU:CD2	1:H:80:LEU:HD12	2.01	0.69
1:G:77:LEU:HD21	1:G:80:LEU:CD1	2.20	0.69
1:A:174:GLU:CG	1:A:234:ARG:N	2.55	0.69
1:B:174:GLU:CG	1:B:234:ARG:N	2.55	0.69
1:I:174:GLU:CG	1:I:234:ARG:N	2.55	0.69
1:D:174:GLU:CG	1:D:234:ARG:N	2.55	0.69
1:J:174:GLU:HG2	1:J:234:ARG:HG3	1.75	0.69
1:K:174:GLU:HG2	1:K:234:ARG:HG3	1.75	0.68
1:B:174:GLU:HG2	1:B:234:ARG:HG3	1.75	0.68
1:H:174:GLU:HG2	1:H:234:ARG:HG3	1.75	0.68
1:G:174:GLU:HG2	1:G:234:ARG:HG3	1.75	0.67
1:I:174:GLU:HG2	1:I:234:ARG:HG3	1.75	0.67
1:E:174:GLU:HG2	1:E:234:ARG:HG3	1.75	0.67
1:C:174:GLU:HG2	1:C:234:ARG:HG3	1.75	0.67
1:A:174:GLU:HG2	1:A:234:ARG:HG3	1.75	0.67
1:B:395:PRO:HG3	1:H:81:ARG:HD2	1.76	0.67
1:C:395:PRO:HG3	1:I:81:ARG:HD2	1.75	0.67
1:D:174:GLU:HG2	1:D:234:ARG:HG3	1.75	0.67
1:F:174:GLU:HG2	1:F:234:ARG:HG3	1.75	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:108:ASP:O	1:K:112:THR:HG23	1.97	0.65
1:E:108:ASP:O	1:E:112:THR:HG23	1.97	0.65
1:F:108:ASP:O	1:F:112:THR:HG23	1.97	0.65
1:A:108:ASP:O	1:A:112:THR:HG23	1.97	0.65
1:D:108:ASP:O	1:D:112:THR:HG23	1.97	0.65
1:I:108:ASP:O	1:I:112:THR:HG23	1.97	0.65
1:J:108:ASP:O	1:J:112:THR:HG23	1.97	0.65
1:H:108:ASP:O	1:H:112:THR:HG23	1.97	0.64
1:B:108:ASP:O	1:B:112:THR:HG23	1.97	0.64
1:I:77:LEU:CD2	1:I:80:LEU:HD12	2.01	0.64
1:G:108:ASP:O	1:G:112:THR:HG23	1.97	0.64
1:C:108:ASP:O	1:C:112:THR:HG23	1.97	0.63
1:H:77:LEU:CD2	1:H:80:LEU:HD11	2.29	0.63
1:F:174:GLU:HG3	1:F:233:VAL:CA	2.29	0.63
1:B:174:GLU:HG3	1:B:233:VAL:CA	2.29	0.63
1:D:77:LEU:CD2	1:D:80:LEU:HD11	2.29	0.63
1:I:174:GLU:HG3	1:I:233:VAL:CA	2.29	0.63
1:J:174:GLU:HG3	1:J:233:VAL:CA	2.29	0.62
1:A:174:GLU:HG3	1:A:233:VAL:CA	2.29	0.62
1:C:77:LEU:CD2	1:C:80:LEU:HD11	2.29	0.62
1:C:174:GLU:HG3	1:C:233:VAL:CA	2.29	0.62
1:G:174:GLU:HG3	1:G:233:VAL:CA	2.29	0.62
1:E:174:GLU:HG3	1:E:233:VAL:CA	2.29	0.62
1:A:77:LEU:CD2	1:A:80:LEU:HD12	2.01	0.62
1:H:174:GLU:HG3	1:H:233:VAL:CA	2.29	0.61
1:D:174:GLU:HG3	1:D:233:VAL:CA	2.29	0.61
1:F:77:LEU:CD2	1:F:80:LEU:HD11	2.29	0.61
1:J:77:LEU:CD2	1:J:80:LEU:HD11	2.29	0.61
1:K:174:GLU:HG3	1:K:233:VAL:CA	2.29	0.61
1:E:77:LEU:CD2	1:E:80:LEU:HD11	2.29	0.60
1:A:77:LEU:CD2	1:A:80:LEU:HD11	2.29	0.60
1:B:229:ILE:HD11	1:B:259:ILE:HD12	1.84	0.60
1:G:77:LEU:CD2	1:G:80:LEU:HD11	2.29	0.60
1:H:229:ILE:HD11	1:H:259:ILE:HD12	1.84	0.60
1:K:229:ILE:HD11	1:K:259:ILE:HD12	1.84	0.60
1:C:229:ILE:HD11	1:C:259:ILE:HD12	1.84	0.60
1:E:229:ILE:HD11	1:E:259:ILE:HD12	1.84	0.60
1:J:229:ILE:HD11	1:J:259:ILE:HD12	1.84	0.60
1:F:229:ILE:HD11	1:F:259:ILE:HD12	1.84	0.60
1:B:174:GLU:HG2	1:B:234:ARG:CG	2.32	0.60
1:D:229:ILE:HD11	1:D:259:ILE:HD12	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:229:ILE:HD11	1:G:259:ILE:HD12	1.84	0.60
1:I:229:ILE:HD11	1:I:259:ILE:HD12	1.84	0.60
1:J:77:LEU:CD2	1:J:80:LEU:HD12	2.01	0.60
1:A:229:ILE:HD11	1:A:259:ILE:HD12	1.84	0.60
1:K:77:LEU:CD2	1:K:80:LEU:HD12	2.01	0.59
1:C:174:GLU:HG2	1:C:234:ARG:CG	2.32	0.59
1:C:331:GLU:OE2	1:C:333:ALA:HB3	2.03	0.59
1:F:174:GLU:HG2	1:F:234:ARG:CG	2.32	0.59
1:C:226:GLY:O	1:C:295:ARG:NH1	2.36	0.59
1:G:174:GLU:HG2	1:G:234:ARG:CG	2.32	0.59
1:I:331:GLU:OE2	1:I:333:ALA:HB3	2.03	0.59
1:F:226:GLY:O	1:F:295:ARG:NH1	2.36	0.59
1:G:331:GLU:OE2	1:G:333:ALA:HB3	2.03	0.59
1:H:331:GLU:OE2	1:H:333:ALA:HB3	2.03	0.59
1:I:174:GLU:HG2	1:I:234:ARG:CG	2.32	0.59
1:I:226:GLY:O	1:I:295:ARG:NH1	2.36	0.59
1:A:331:GLU:OE2	1:A:333:ALA:HB3	2.03	0.59
1:B:331:GLU:OE2	1:B:333:ALA:HB3	2.03	0.59
1:H:226:GLY:O	1:H:295:ARG:NH1	2.36	0.59
1:D:331:GLU:OE2	1:D:333:ALA:HB3	2.03	0.59
1:I:77:LEU:CD2	1:I:80:LEU:HD11	2.29	0.59
1:D:226:GLY:O	1:D:295:ARG:NH1	2.36	0.59
1:K:331:GLU:OE2	1:K:333:ALA:HB3	2.03	0.59
1:F:331:GLU:OE2	1:F:333:ALA:HB3	2.03	0.59
1:B:226:GLY:O	1:B:295:ARG:NH1	2.36	0.59
1:J:174:GLU:HG2	1:J:234:ARG:CG	2.32	0.59
1:A:174:GLU:HG2	1:A:234:ARG:CG	2.32	0.59
1:B:77:LEU:CD2	1:B:80:LEU:HD11	2.28	0.58
1:E:174:GLU:HG2	1:E:234:ARG:CG	2.32	0.58
1:K:226:GLY:O	1:K:295:ARG:NH1	2.36	0.58
1:G:226:GLY:O	1:G:295:ARG:NH1	2.36	0.58
1:H:174:GLU:HG2	1:H:234:ARG:CG	2.32	0.58
1:K:174:GLU:HG2	1:K:234:ARG:CG	2.32	0.58
1:A:305:GLU:OE2	1:A:305:GLU:N	2.29	0.58
1:J:226:GLY:O	1:J:295:ARG:NH1	2.36	0.58
1:J:331:GLU:OE2	1:J:333:ALA:HB3	2.03	0.58
1:D:174:GLU:HG2	1:D:234:ARG:CG	2.32	0.58
1:E:226:GLY:O	1:E:295:ARG:NH1	2.36	0.58
1:A:226:GLY:O	1:A:295:ARG:NH1	2.36	0.58
1:G:305:GLU:OE2	1:G:305:GLU:N	2.29	0.58
1:E:331:GLU:OE2	1:E:333:ALA:HB3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:GLU:OE2	1:B:305:GLU:N	2.29	0.57
1:E:174:GLU:HG3	1:E:234:ARG:N	2.19	0.57
1:E:305:GLU:OE2	1:E:305:GLU:N	2.29	0.57
1:I:174:GLU:HG3	1:I:234:ARG:N	2.20	0.56
1:F:174:GLU:HG3	1:F:234:ARG:N	2.20	0.56
1:A:174:GLU:HG3	1:A:234:ARG:N	2.20	0.56
1:B:174:GLU:HG3	1:B:234:ARG:N	2.19	0.56
1:H:305:GLU:OE2	1:H:305:GLU:N	2.29	0.56
1:K:77:LEU:CD2	1:K:80:LEU:HD11	2.29	0.56
1:B:382:VAL:HG13	1:B:382:VAL:O	2.06	0.56
1:G:382:VAL:O	1:G:382:VAL:HG13	2.06	0.56
1:A:382:VAL:O	1:A:382:VAL:HG13	2.06	0.56
1:F:382:VAL:O	1:F:382:VAL:HG13	2.06	0.55
1:H:174:GLU:HG3	1:H:234:ARG:N	2.20	0.55
1:H:382:VAL:O	1:H:382:VAL:HG13	2.06	0.55
1:J:174:GLU:HG3	1:J:234:ARG:N	2.20	0.55
1:D:382:VAL:HG13	1:D:382:VAL:O	2.06	0.55
1:E:382:VAL:HG13	1:E:382:VAL:O	2.06	0.55
1:G:174:GLU:HG3	1:G:234:ARG:N	2.20	0.55
1:I:382:VAL:O	1:I:382:VAL:HG13	2.06	0.55
1:K:382:VAL:HG13	1:K:382:VAL:O	2.06	0.55
1:C:305:GLU:OE2	1:C:305:GLU:N	2.29	0.55
1:J:382:VAL:HG13	1:J:382:VAL:O	2.06	0.55
1:C:174:GLU:HG3	1:C:234:ARG:N	2.20	0.55
1:K:174:GLU:HG3	1:K:234:ARG:N	2.20	0.55
1:D:293:HIS:ND1	1:I:164:TYR:OH	2.40	0.54
1:C:382:VAL:HG13	1:C:382:VAL:O	2.06	0.54
1:G:181:LEU:HD23	1:G:182:HIS:N	2.23	0.54
1:F:181:LEU:HD23	1:F:182:HIS:N	2.23	0.54
1:K:181:LEU:HD23	1:K:182:HIS:N	2.23	0.54
1:C:293:HIS:ND1	1:H:164:TYR:OH	2.41	0.53
1:F:195:ASP:OD1	1:A:320:ASN:ND2	2.41	0.53
1:K:305:GLU:OE2	1:K:305:GLU:N	2.29	0.53
1:I:181:LEU:HD23	1:I:182:HIS:N	2.23	0.53
1:D:181:LEU:HD23	1:D:182:HIS:N	2.23	0.53
1:H:181:LEU:HD23	1:H:182:HIS:N	2.23	0.53
1:C:181:LEU:HD23	1:C:182:HIS:N	2.23	0.53
1:E:181:LEU:HD23	1:E:182:HIS:N	2.23	0.53
1:A:181:LEU:HD23	1:A:182:HIS:N	2.23	0.53
1:I:305:GLU:OE2	1:I:305:GLU:N	2.29	0.53
1:J:181:LEU:HD23	1:J:182:HIS:N	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:LEU:HD23	1:B:182:HIS:N	2.23	0.53
1:C:320:ASN:ND2	1:H:195:ASP:OD1	2.42	0.52
1:D:320:ASN:ND2	1:I:195:ASP:OD1	2.42	0.52
1:S:775:TYR:O	1:S:779:ASP:N	2.34	0.52
1:D:262:CYS:SG	1:I:394:MET:HE1	2.50	0.52
1:U:775:TYR:O	1:U:779:ASP:N	2.34	0.52
1:B:293:HIS:ND1	1:G:164:TYR:OH	2.42	0.52
1:D:174:GLU:HG3	1:D:234:ARG:N	2.20	0.52
1:H:174:GLU:HG2	1:H:234:ARG:N	2.25	0.52
1:C:262:CYS:SG	1:H:394:MET:HE1	2.50	0.51
1:D:305:GLU:OE2	1:D:305:GLU:N	2.29	0.51
1:R:775:TYR:O	1:R:779:ASP:N	2.34	0.51
1:V:775:TYR:O	1:V:779:ASP:N	2.34	0.51
1:D:174:GLU:HG2	1:D:234:ARG:N	2.25	0.51
1:C:174:GLU:HG2	1:C:234:ARG:N	2.25	0.51
1:K:174:GLU:HG2	1:K:234:ARG:N	2.25	0.51
1:J:174:GLU:HG2	1:J:234:ARG:N	2.25	0.51
1:A:174:GLU:HG2	1:A:234:ARG:N	2.25	0.51
1:B:174:GLU:HG2	1:B:234:ARG:N	2.25	0.51
1:L:458:LYS:H	1:L:458:LYS:HD2	1.76	0.51
1:J:174:GLU:CG	1:J:233:VAL:C	2.82	0.51
1:V:458:LYS:H	1:V:458:LYS:HD2	1.76	0.51
1:L:775:TYR:O	1:L:779:ASP:N	2.34	0.50
1:T:458:LYS:H	1:T:458:LYS:HD2	1.76	0.50
1:H:174:GLU:CG	1:H:233:VAL:C	2.82	0.50
1:I:174:GLU:HG2	1:I:234:ARG:N	2.25	0.50
1:M:458:LYS:H	1:M:458:LYS:HD2	1.76	0.50
1:R:458:LYS:H	1:R:458:LYS:HD2	1.76	0.50
1:S:458:LYS:H	1:S:458:LYS:HD2	1.76	0.50
1:U:458:LYS:H	1:U:458:LYS:HD2	1.76	0.50
1:Q:775:TYR:O	1:Q:779:ASP:N	2.34	0.50
1:B:320:ASN:ND2	1:G:195:ASP:OD1	2.43	0.50
1:C:174:GLU:CG	1:C:233:VAL:C	2.82	0.50
1:F:305:GLU:OE2	1:F:305:GLU:N	2.29	0.50
1:Q:458:LYS:H	1:Q:458:LYS:HD2	1.76	0.50
1:F:174:GLU:HG2	1:F:234:ARG:N	2.25	0.50
1:N:458:LYS:H	1:N:458:LYS:HD2	1.76	0.50
1:P:458:LYS:H	1:P:458:LYS:HD2	1.76	0.50
1:M:775:TYR:O	1:M:779:ASP:N	2.34	0.49
1:G:174:GLU:HG2	1:G:234:ARG:N	2.25	0.49
1:D:164:TYR:OH	1:J:293:HIS:ND1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:77:LEU:CD2	1:F:80:LEU:HD12	2.01	0.49
1:P:775:TYR:O	1:P:779:ASP:N	2.34	0.49
1:N:775:TYR:O	1:N:779:ASP:N	2.34	0.49
1:E:174:GLU:HG2	1:E:234:ARG:N	2.25	0.49
1:I:174:GLU:HG3	1:I:233:VAL:HA	1.94	0.49
1:J:174:GLU:HG3	1:J:233:VAL:HA	1.94	0.49
1:C:174:GLU:HG3	1:C:233:VAL:HA	1.94	0.49
1:O:458:LYS:H	1:O:458:LYS:HD2	1.76	0.49
1:E:195:ASP:OD1	1:K:320:ASN:ND2	2.44	0.49
1:D:174:GLU:HG3	1:D:233:VAL:HA	1.94	0.49
1:E:174:GLU:CG	1:E:233:VAL:C	2.82	0.48
1:J:305:GLU:OE2	1:J:305:GLU:N	2.29	0.48
1:O:775:TYR:O	1:O:779:ASP:N	2.34	0.48
1:B:262:CYS:SG	1:G:394:MET:HE1	2.53	0.48
1:E:320:ASN:ND2	1:J:195:ASP:OD1	2.44	0.48
1:F:164:TYR:OH	1:A:293:HIS:ND1	2.46	0.48
1:F:174:GLU:HG3	1:F:233:VAL:HA	1.94	0.48
1:A:174:GLU:HG3	1:A:233:VAL:HA	1.95	0.48
1:E:293:HIS:ND1	1:J:164:TYR:OH	2.47	0.48
1:K:174:GLU:HG3	1:K:233:VAL:HA	1.94	0.48
1:T:775:TYR:O	1:T:779:ASP:N	2.34	0.48
1:B:174:GLU:HG3	1:B:233:VAL:HA	1.94	0.48
1:G:293:HIS:ND1	1:A:164:TYR:OH	2.46	0.48
1:F:174:GLU:CG	1:F:233:VAL:C	2.82	0.48
1:F:293:HIS:ND1	1:K:164:TYR:OH	2.47	0.47
1:G:368:TYR:CE1	1:G:372:ILE:HD11	2.49	0.47
1:H:174:GLU:HG3	1:H:233:VAL:HA	1.95	0.47
1:A:368:TYR:CE1	1:A:372:ILE:HD11	2.49	0.47
1:G:174:GLU:HG3	1:G:233:VAL:HA	1.94	0.47
1:E:174:GLU:HG3	1:E:233:VAL:HA	1.94	0.47
1:B:368:TYR:CE1	1:B:372:ILE:HD11	2.49	0.47
1:C:195:ASP:OD1	1:I:320:ASN:ND2	2.47	0.47
1:C:368:TYR:CE1	1:C:372:ILE:HD11	2.49	0.47
1:D:228:PHE:O	1:D:229:ILE:HD13	2.15	0.47
1:F:368:TYR:CE1	1:F:372:ILE:HD11	2.49	0.47
1:H:228:PHE:O	1:H:229:ILE:HD13	2.15	0.47
1:I:228:PHE:O	1:I:229:ILE:HD13	2.15	0.47
1:I:368:TYR:CE1	1:I:372:ILE:HD11	2.49	0.47
1:J:228:PHE:O	1:J:229:ILE:HD13	2.15	0.47
1:A:174:GLU:CG	1:A:233:VAL:C	2.82	0.47
1:B:72:THR:O	1:B:72:THR:HG22	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:368:TYR:CE1	1:H:372:ILE:HD11	2.49	0.47
1:C:228:PHE:O	1:C:229:ILE:HD13	2.15	0.47
1:J:368:TYR:CE1	1:J:372:ILE:HD11	2.49	0.47
1:D:368:TYR:CE1	1:D:372:ILE:HD11	2.49	0.47
1:K:228:PHE:O	1:K:229:ILE:HD13	2.15	0.47
1:D:394:MET:HE1	1:J:262:CYS:SG	2.55	0.46
1:E:368:TYR:CE1	1:E:372:ILE:HD11	2.49	0.46
1:F:72:THR:HG22	1:F:72:THR:O	2.15	0.46
1:G:228:PHE:O	1:G:229:ILE:HD13	2.15	0.46
1:H:72:THR:HG22	1:H:72:THR:O	2.15	0.46
1:E:228:PHE:O	1:E:229:ILE:HD13	2.15	0.46
1:K:368:TYR:CE1	1:K:372:ILE:HD11	2.49	0.46
1:K:229:ILE:HD11	1:K:259:ILE:CD1	2.46	0.46
1:E:229:ILE:HD11	1:E:259:ILE:CD1	2.46	0.46
1:F:229:ILE:HD11	1:F:259:ILE:CD1	2.46	0.46
1:J:229:ILE:HD11	1:J:259:ILE:CD1	2.46	0.46
1:G:72:THR:O	1:G:72:THR:HG22	2.15	0.46
1:J:72:THR:HG22	1:J:72:THR:O	2.15	0.46
1:S:458:LYS:HD2	1:S:458:LYS:N	2.31	0.46
1:C:229:ILE:HD11	1:C:259:ILE:CD1	2.46	0.46
1:D:229:ILE:HD11	1:D:259:ILE:CD1	2.46	0.46
1:F:228:PHE:O	1:F:229:ILE:HD13	2.15	0.46
1:H:229:ILE:HD11	1:H:259:ILE:CD1	2.46	0.46
1:I:229:ILE:HD11	1:I:259:ILE:CD1	2.46	0.46
1:A:229:ILE:HD11	1:A:259:ILE:CD1	2.46	0.46
1:M:458:LYS:HD2	1:M:458:LYS:N	2.31	0.46
1:R:458:LYS:HD2	1:R:458:LYS:N	2.31	0.46
1:T:458:LYS:HD2	1:T:458:LYS:N	2.31	0.46
1:B:229:ILE:HD11	1:B:259:ILE:CD1	2.46	0.46
1:G:174:GLU:CG	1:G:233:VAL:C	2.82	0.46
1:G:229:ILE:HD11	1:G:259:ILE:CD1	2.46	0.46
1:K:174:GLU:CG	1:K:233:VAL:C	2.82	0.46
1:A:228:PHE:O	1:A:229:ILE:HD13	2.15	0.46
1:N:458:LYS:HD2	1:N:458:LYS:N	2.31	0.46
1:C:72:THR:HG22	1:C:72:THR:O	2.15	0.46
1:G:262:CYS:SG	1:A:394:MET:HE1	2.55	0.46
1:B:228:PHE:O	1:B:229:ILE:HD13	2.15	0.46
1:B:339:LYS:O	1:B:343:ASP:OD1	2.34	0.46
1:D:195:ASP:OD1	1:J:320:ASN:ND2	2.48	0.46
1:Q:458:LYS:HD2	1:Q:458:LYS:N	2.31	0.46
1:B:195:ASP:OD1	1:H:320:ASN:ND2	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:339:LYS:O	1:H:343:ASP:OD1	2.34	0.45
1:K:72:THR:HG22	1:K:72:THR:O	2.15	0.45
1:L:458:LYS:HD2	1:L:458:LYS:N	2.31	0.45
1:U:458:LYS:HD2	1:U:458:LYS:N	2.31	0.45
1:C:339:LYS:O	1:C:343:ASP:OD1	2.34	0.45
1:E:339:LYS:O	1:E:343:ASP:OD1	2.34	0.45
1:G:339:LYS:O	1:G:343:ASP:OD1	2.34	0.45
1:J:339:LYS:O	1:J:343:ASP:OD1	2.35	0.45
1:D:339:LYS:O	1:D:343:ASP:OD1	2.34	0.45
1:F:339:LYS:O	1:F:343:ASP:OD1	2.34	0.45
1:G:331:GLU:OE2	1:G:333:ALA:N	2.50	0.45
1:J:331:GLU:OE2	1:J:333:ALA:N	2.50	0.45
1:K:331:GLU:OE2	1:K:333:ALA:N	2.50	0.45
1:A:72:THR:HG22	1:A:72:THR:O	2.15	0.45
1:E:243:GLN:O	1:E:247:ARG:N	2.45	0.45
1:I:331:GLU:OE2	1:I:333:ALA:N	2.50	0.45
1:I:339:LYS:O	1:I:343:ASP:OD1	2.34	0.45
1:K:339:LYS:O	1:K:343:ASP:OD1	2.35	0.45
1:C:331:GLU:OE2	1:C:333:ALA:N	2.50	0.45
1:G:320:ASN:ND2	1:A:195:ASP:OD1	2.48	0.45
1:J:243:GLN:O	1:J:247:ARG:N	2.45	0.45
1:A:331:GLU:OE2	1:A:333:ALA:N	2.50	0.45
1:A:339:LYS:O	1:A:343:ASP:OD1	2.34	0.45
1:C:322:LEU:HD23	1:C:339:LYS:HD3	1.98	0.45
1:E:72:THR:HG22	1:E:72:THR:O	2.15	0.45
1:P:458:LYS:HD2	1:P:458:LYS:N	2.31	0.45
1:E:262:CYS:SG	1:J:394:MET:HE1	2.57	0.45
1:H:322:LEU:HD23	1:H:339:LYS:HD3	1.98	0.45
1:K:174:GLU:CD	1:K:234:ARG:HA	2.24	0.45
1:O:458:LYS:HD2	1:O:458:LYS:N	2.31	0.45
1:V:458:LYS:HD2	1:V:458:LYS:N	2.31	0.45
1:F:394:MET:HE1	1:A:262:CYS:SG	2.56	0.45
1:B:331:GLU:OE2	1:B:333:ALA:N	2.50	0.45
1:D:72:THR:O	1:D:72:THR:HG22	2.15	0.45
1:I:72:THR:O	1:I:72:THR:HG22	2.15	0.45
1:K:243:GLN:O	1:K:247:ARG:N	2.45	0.45
1:D:331:GLU:OE2	1:D:333:ALA:N	2.50	0.45
1:F:331:GLU:OE2	1:F:333:ALA:N	2.50	0.44
1:B:322:LEU:HD23	1:B:339:LYS:HD3	1.98	0.44
1:C:162:VAL:O	1:C:166:VAL:HG22	2.18	0.44
1:F:174:GLU:CD	1:F:234:ARG:HA	2.24	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:331:GLU:OE2	1:H:333:ALA:N	2.50	0.44
1:I:322:LEU:HD23	1:I:339:LYS:HD3	1.98	0.44
1:J:162:VAL:O	1:J:166:VAL:HG22	2.18	0.44
1:B:162:VAL:O	1:B:166:VAL:HG22	2.18	0.44
1:D:243:GLN:O	1:D:247:ARG:N	2.45	0.44
1:E:174:GLU:CD	1:E:234:ARG:HA	2.24	0.44
1:E:331:GLU:OE2	1:E:333:ALA:N	2.50	0.44
1:F:322:LEU:HD23	1:F:339:LYS:HD3	1.98	0.44
1:K:162:VAL:O	1:K:166:VAL:HG22	2.18	0.44
1:A:322:LEU:HD23	1:A:339:LYS:HD3	1.99	0.44
1:C:228:PHE:C	1:C:229:ILE:HD13	2.43	0.44
1:F:228:PHE:C	1:F:229:ILE:HD13	2.43	0.44
1:G:162:VAL:O	1:G:166:VAL:HG22	2.18	0.44
1:I:243:GLN:O	1:I:247:ARG:N	2.45	0.44
1:K:322:LEU:HD23	1:K:339:LYS:HD3	1.98	0.44
1:E:322:LEU:HD23	1:E:339:LYS:HD3	1.98	0.44
1:J:322:LEU:HD23	1:J:339:LYS:HD3	1.98	0.44
1:A:162:VAL:O	1:A:166:VAL:HG22	2.18	0.44
1:D:162:VAL:O	1:D:166:VAL:HG22	2.18	0.44
1:E:164:TYR:OH	1:K:293:HIS:ND1	2.51	0.44
1:F:243:GLN:O	1:F:247:ARG:N	2.45	0.44
1:G:322:LEU:HD23	1:G:339:LYS:HD3	1.98	0.44
1:D:322:LEU:HD23	1:D:339:LYS:HD3	1.98	0.43
1:H:110:SER:O	1:H:114:VAL:HG22	2.18	0.43
1:F:162:VAL:O	1:F:166:VAL:HG22	2.18	0.43
1:I:228:PHE:C	1:I:229:ILE:HD13	2.43	0.43
1:B:228:PHE:C	1:B:229:ILE:HD13	2.43	0.43
1:E:181:LEU:HD22	1:E:183:THR:HB	2.00	0.43
1:J:110:SER:O	1:J:114:VAL:HG22	2.19	0.43
1:J:228:PHE:C	1:J:229:ILE:HD13	2.43	0.43
1:B:110:SER:O	1:B:114:VAL:HG22	2.18	0.43
1:C:110:SER:O	1:C:114:VAL:HG22	2.18	0.43
1:G:228:PHE:C	1:G:229:ILE:HD13	2.43	0.43
1:I:162:VAL:O	1:I:166:VAL:HG22	2.18	0.43
1:E:228:PHE:C	1:E:229:ILE:HD13	2.43	0.43
1:F:262:CYS:SG	1:K:394:MET:HE1	2.58	0.43
1:H:162:VAL:O	1:H:166:VAL:HG22	2.18	0.43
1:H:228:PHE:C	1:H:229:ILE:HD13	2.43	0.43
1:K:181:LEU:HD22	1:K:183:THR:HB	2.00	0.43
1:A:228:PHE:C	1:A:229:ILE:HD13	2.43	0.43
1:F:110:SER:O	1:F:114:VAL:HG22	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:110:SER:O	1:I:114:VAL:HG22	2.18	0.43
1:K:228:PHE:C	1:K:229:ILE:HD13	2.43	0.43
1:A:110:SER:O	1:A:114:VAL:HG22	2.19	0.43
1:L:465:LYS:HE3	1:L:465:LYS:HB3	1.90	0.43
1:F:181:LEU:HD22	1:F:183:THR:HB	2.00	0.43
1:J:181:LEU:HD22	1:J:183:THR:HB	2.01	0.43
1:E:162:VAL:O	1:E:166:VAL:HG22	2.18	0.43
1:A:181:LEU:HD22	1:A:183:THR:HB	2.00	0.43
1:B:164:TYR:OH	1:H:293:HIS:ND1	2.52	0.43
1:D:110:SER:O	1:D:114:VAL:HG22	2.19	0.43
1:E:110:SER:O	1:E:114:VAL:HG22	2.19	0.43
1:U:465:LYS:HE3	1:U:465:LYS:HB3	1.90	0.43
1:C:243:GLN:O	1:C:247:ARG:N	2.45	0.43
1:A:77:LEU:HD21	1:A:80:LEU:HD11	1.96	0.43
1:A:174:GLU:CD	1:A:234:ARG:HA	2.24	0.43
1:G:110:SER:O	1:G:114:VAL:HG22	2.18	0.42
1:I:111:ILE:O	1:I:114:VAL:CG2	2.67	0.42
1:I:174:GLU:CG	1:I:233:VAL:C	2.82	0.42
1:J:174:GLU:CD	1:J:234:ARG:HA	2.24	0.42
1:A:221:ASP:OD1	1:A:224:GLU:HB2	2.20	0.42
1:D:181:LEU:HD22	1:D:183:THR:HB	2.00	0.42
1:D:228:PHE:C	1:D:229:ILE:HD13	2.43	0.42
1:E:111:ILE:O	1:E:114:VAL:CG2	2.67	0.42
1:E:394:MET:HE1	1:K:262:CYS:SG	2.59	0.42
1:F:111:ILE:O	1:F:114:VAL:CG2	2.67	0.42
1:A:243:GLN:O	1:A:247:ARG:N	2.45	0.42
1:B:111:ILE:O	1:B:114:VAL:CG2	2.67	0.42
1:B:181:LEU:HD22	1:B:183:THR:HB	2.00	0.42
1:F:77:LEU:HD21	1:F:80:LEU:HD11	1.96	0.42
1:F:320:ASN:ND2	1:K:195:ASP:OD1	2.49	0.42
1:H:111:ILE:O	1:H:114:VAL:CG2	2.67	0.42
1:J:111:ILE:O	1:J:114:VAL:CG2	2.67	0.42
1:J:176:MET:HG3	1:J:176:MET:O	2.20	0.42
1:B:221:ASP:OD1	1:B:224:GLU:HB2	2.20	0.42
1:C:164:TYR:OH	1:I:293:HIS:ND1	2.52	0.42
1:C:181:LEU:HD22	1:C:183:THR:HB	2.00	0.42
1:E:176:MET:HG3	1:E:176:MET:O	2.20	0.42
1:G:181:LEU:HD22	1:G:183:THR:HB	2.00	0.42
1:H:221:ASP:OD1	1:H:224:GLU:HB2	2.19	0.42
1:C:176:MET:O	1:C:176:MET:HG3	2.20	0.42
1:D:176:MET:O	1:D:176:MET:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:77:LEU:HD21	1:G:80:LEU:HD11	1.96	0.42
1:K:111:ILE:O	1:K:114:VAL:CG2	2.67	0.42
1:A:111:ILE:O	1:A:114:VAL:CG2	2.67	0.42
1:B:176:MET:HG3	1:B:176:MET:O	2.20	0.42
1:C:111:ILE:O	1:C:114:VAL:CG2	2.67	0.42
1:G:111:ILE:O	1:G:114:VAL:CG2	2.67	0.42
1:H:176:MET:O	1:H:176:MET:HG3	2.20	0.42
1:I:176:MET:HG3	1:I:176:MET:O	2.20	0.42
1:I:181:LEU:HD22	1:I:183:THR:HB	2.00	0.42
1:K:110:SER:O	1:K:114:VAL:HG22	2.19	0.42
1:K:176:MET:O	1:K:176:MET:HG3	2.20	0.42
1:F:176:MET:O	1:F:176:MET:HG3	2.20	0.42
1:H:189:VAL:O	1:H:190:SER:OG	2.37	0.42
1:G:176:MET:HG3	1:G:176:MET:O	2.20	0.42
1:A:176:MET:O	1:A:176:MET:HG3	2.20	0.42
1:E:221:ASP:OD1	1:E:224:GLU:HB2	2.20	0.42
1:F:221:ASP:OD1	1:F:224:GLU:HB2	2.20	0.42
1:H:243:GLN:O	1:H:247:ARG:N	2.45	0.42
1:K:77:LEU:HD21	1:K:80:LEU:HD11	1.96	0.42
1:K:385:ILE:HG23	1:K:385:ILE:O	2.20	0.42
1:P:465:LYS:HE3	1:P:465:LYS:HB3	1.90	0.42
1:H:181:LEU:HD22	1:H:183:THR:HB	2.01	0.41
1:K:221:ASP:OD1	1:K:224:GLU:HB2	2.20	0.41
1:D:111:ILE:O	1:D:114:VAL:CG2	2.67	0.41
1:B:189:VAL:O	1:B:190:SER:OG	2.37	0.41
1:D:221:ASP:OD1	1:D:224:GLU:HB2	2.19	0.41
1:G:221:ASP:OD1	1:G:224:GLU:HB2	2.20	0.41
1:G:243:GLN:O	1:G:247:ARG:N	2.45	0.41
1:I:221:ASP:OD1	1:I:224:GLU:HB2	2.19	0.41
1:J:385:ILE:O	1:J:385:ILE:HG23	2.20	0.41
1:B:77:LEU:HD21	1:B:80:LEU:HD11	1.96	0.41
1:A:385:ILE:O	1:A:385:ILE:HG23	2.20	0.41
1:B:243:GLN:O	1:B:247:ARG:N	2.45	0.41
1:J:221:ASP:OD1	1:J:224:GLU:HB2	2.19	0.41
1:F:385:ILE:O	1:F:385:ILE:HG23	2.20	0.41
1:C:385:ILE:O	1:C:385:ILE:HG23	2.20	0.41
1:E:385:ILE:O	1:E:385:ILE:HG23	2.20	0.41
1:F:260:HIS:O	1:F:275:TYR:N	2.51	0.41
1:C:221:ASP:OD1	1:C:224:GLU:HB2	2.19	0.41
1:C:77:LEU:HD21	1:C:80:LEU:HD11	1.96	0.41
1:E:275:TYR:HE2	1:E:277:ILE:HD11	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:275:TYR:HE2	1:F:277:ILE:HD11	1.86	0.41
1:G:174:GLU:CD	1:G:234:ARG:HA	2.24	0.41
1:A:275:TYR:HE2	1:A:277:ILE:HD11	1.86	0.41
1:F:111:ILE:O	1:F:114:VAL:HG22	2.22	0.40
1:H:111:ILE:O	1:H:114:VAL:HG22	2.22	0.40
1:I:385:ILE:O	1:I:385:ILE:HG23	2.20	0.40
1:B:111:ILE:O	1:B:114:VAL:HG22	2.22	0.40
1:D:174:GLU:CG	1:D:233:VAL:C	2.82	0.40
1:D:174:GLU:CD	1:D:234:ARG:HA	2.24	0.40
1:G:111:ILE:O	1:G:114:VAL:HG22	2.22	0.40
1:H:385:ILE:HG23	1:H:385:ILE:O	2.20	0.40
1:A:111:ILE:O	1:A:114:VAL:HG22	2.22	0.40
1:A:152:ALA:O	1:A:177:ARG:NH2	2.54	0.40
1:C:316:VAL:HG22	1:C:322:LEU:CD1	2.51	0.40
1:E:77:LEU:HD21	1:E:80:LEU:HD11	1.96	0.40
1:F:316:VAL:HG22	1:F:322:LEU:CD1	2.51	0.40
1:I:152:ALA:O	1:I:177:ARG:NH2	2.54	0.40
1:J:152:ALA:O	1:J:177:ARG:NH2	2.54	0.40
1:J:275:TYR:HE2	1:J:277:ILE:HD11	1.86	0.40
1:B:316:VAL:HG22	1:B:322:LEU:CD1	2.51	0.40
1:B:385:ILE:O	1:B:385:ILE:HG23	2.20	0.40
1:D:152:ALA:O	1:D:177:ARG:NH2	2.54	0.40
1:D:385:ILE:O	1:D:385:ILE:HG23	2.20	0.40
1:H:316:VAL:HG22	1:H:322:LEU:CD1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	320/1011 (32%)	309 (97%)	11 (3%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	320/1011 (32%)	308 (96%)	12 (4%)	0	100	100
1	C	320/1011 (32%)	308 (96%)	12 (4%)	0	100	100
1	D	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	E	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	F	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	G	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	H	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	I	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	J	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	K	320/1011 (32%)	309 (97%)	11 (3%)	0	100	100
1	L	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	M	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	N	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	O	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	P	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	Q	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	R	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	S	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	T	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	U	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
1	V	325/1011 (32%)	319 (98%)	6 (2%)	0	100	100
All	All	7095/22242 (32%)	6906 (97%)	189 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	286/877 (33%)	286 (100%)	0	100	100
1	B	286/877 (33%)	286 (100%)	0	100	100
1	C	286/877 (33%)	286 (100%)	0	100	100
1	D	286/877 (33%)	286 (100%)	0	100	100
1	E	286/877 (33%)	286 (100%)	0	100	100
1	F	286/877 (33%)	286 (100%)	0	100	100
1	G	286/877 (33%)	286 (100%)	0	100	100
1	H	286/877 (33%)	286 (100%)	0	100	100
1	I	286/877 (33%)	286 (100%)	0	100	100
1	J	286/877 (33%)	286 (100%)	0	100	100
1	K	286/877 (33%)	286 (100%)	0	100	100
1	L	17/877 (2%)	17 (100%)	0	100	100
1	M	17/877 (2%)	17 (100%)	0	100	100
1	N	17/877 (2%)	17 (100%)	0	100	100
1	O	17/877 (2%)	17 (100%)	0	100	100
1	P	17/877 (2%)	17 (100%)	0	100	100
1	Q	17/877 (2%)	17 (100%)	0	100	100
1	R	17/877 (2%)	17 (100%)	0	100	100
1	S	17/877 (2%)	17 (100%)	0	100	100
1	T	17/877 (2%)	17 (100%)	0	100	100
1	U	17/877 (2%)	17 (100%)	0	100	100
1	V	17/877 (2%)	17 (100%)	0	100	100
All	All	3333/19294 (17%)	3333 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	93	HIS
1	D	185	ASN
1	H	93	HIS
1	A	93	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

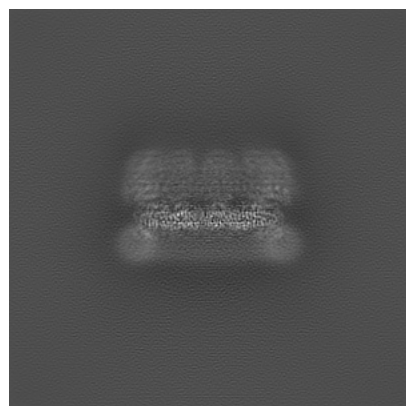
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29290. These allow visual inspection of the internal detail of the map and identification of artifacts.

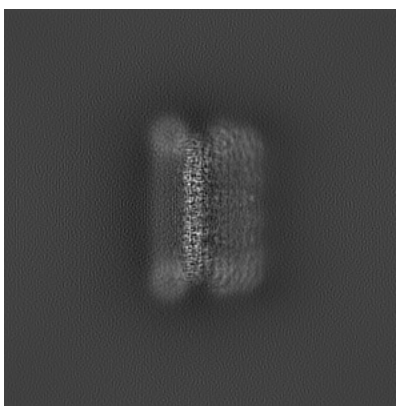
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

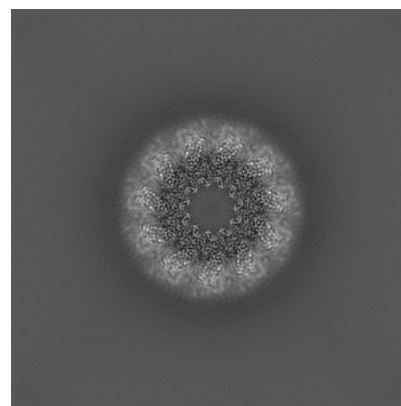
6.1.1 Primary map



X

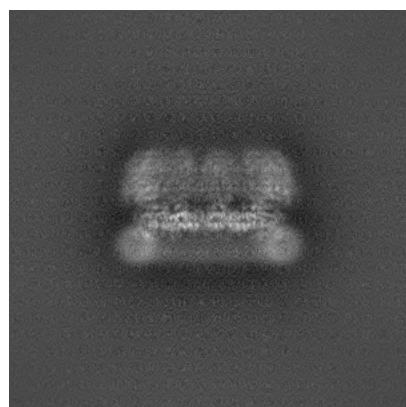


Y

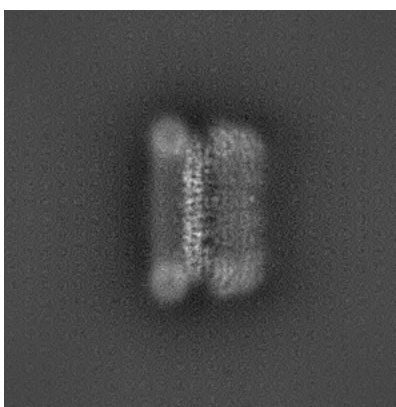


Z

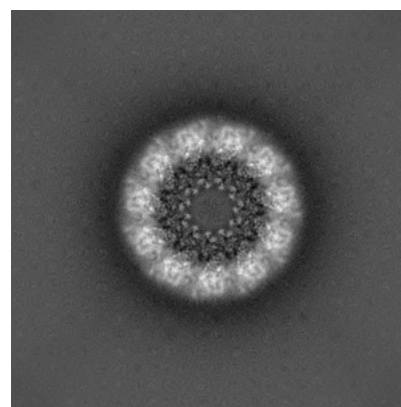
6.1.2 Raw map



X



Y

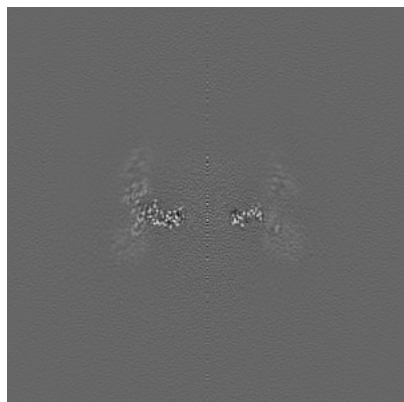


Z

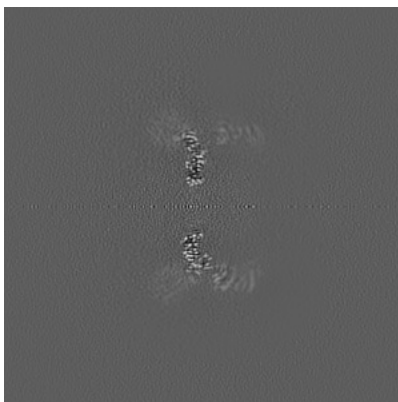
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

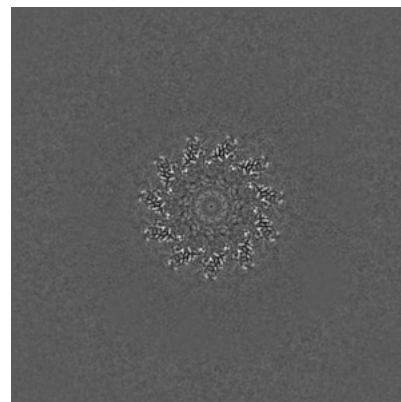
6.2.1 Primary map



X Index: 288



Y Index: 288



Z Index: 288

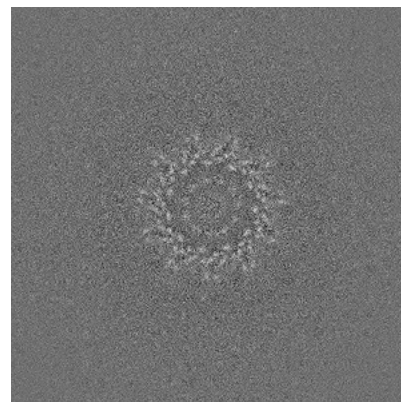
6.2.2 Raw map



X Index: 288



Y Index: 288

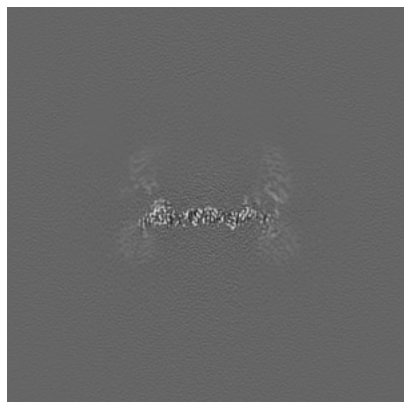


Z Index: 288

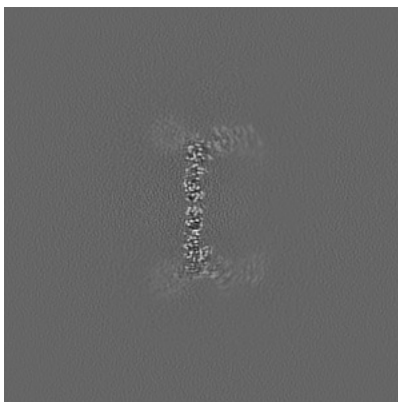
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

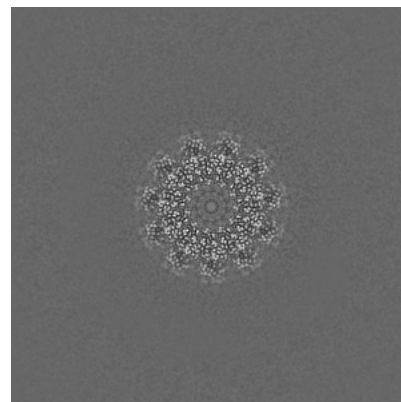
6.3.1 Primary map



X Index: 327

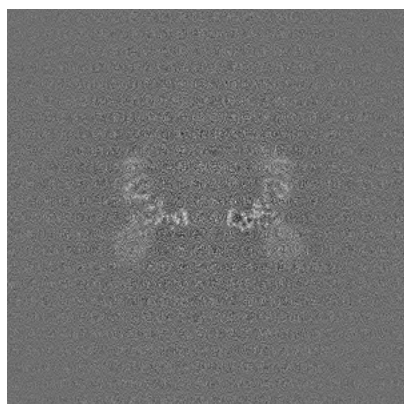


Y Index: 253



Z Index: 277

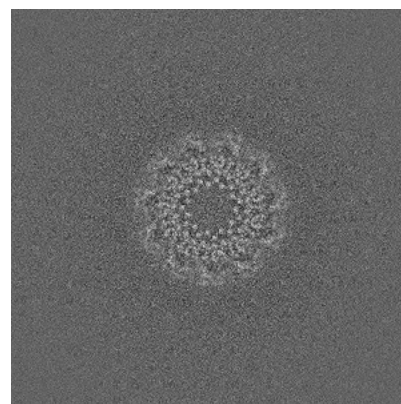
6.3.2 Raw map



X Index: 304



Y Index: 288

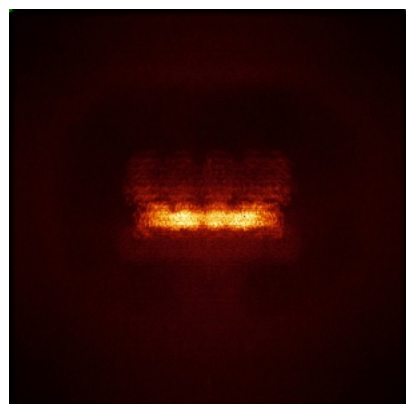


Z Index: 277

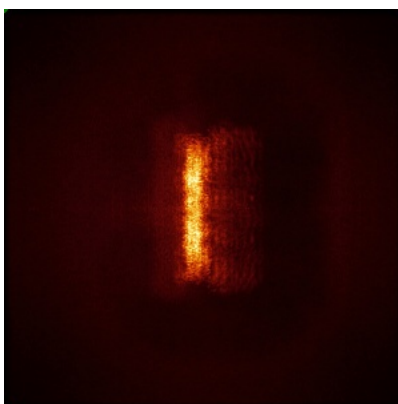
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

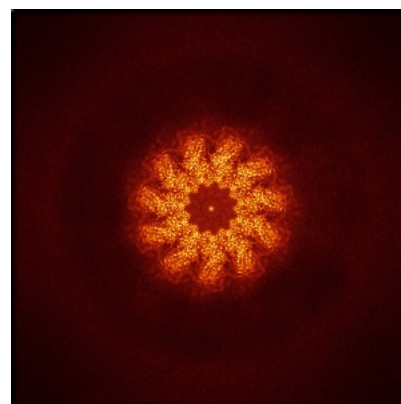
6.4.1 Primary map



X

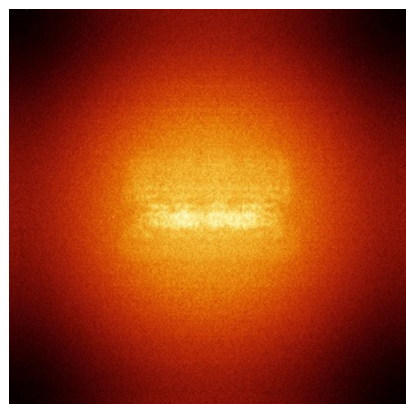


Y

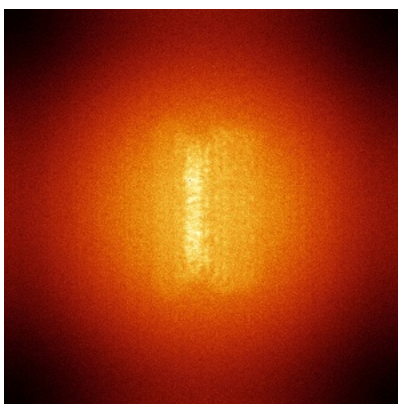


Z

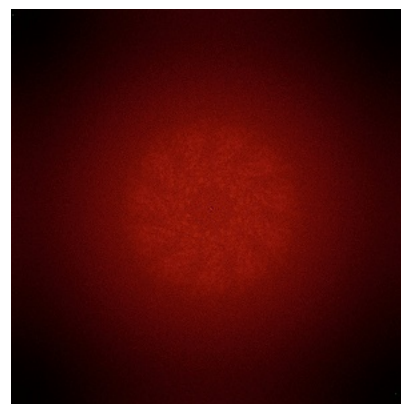
6.4.2 Raw map



X



Y

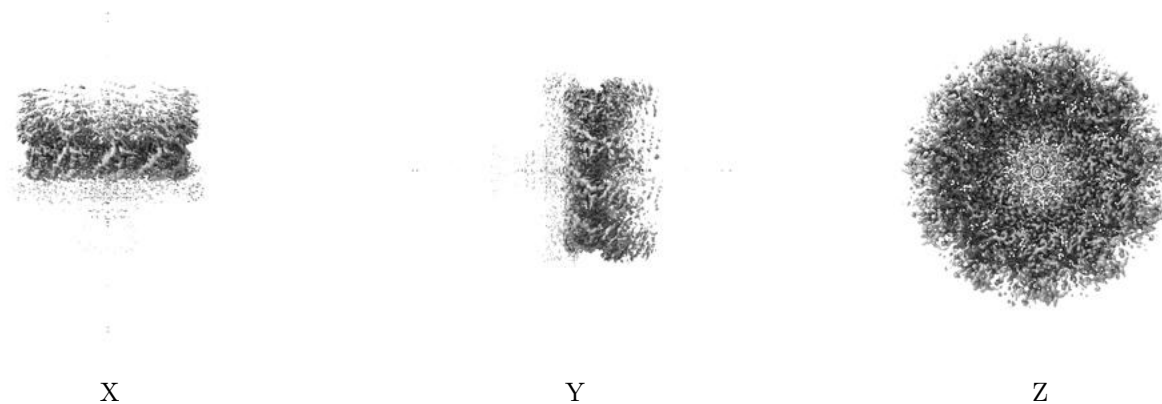


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

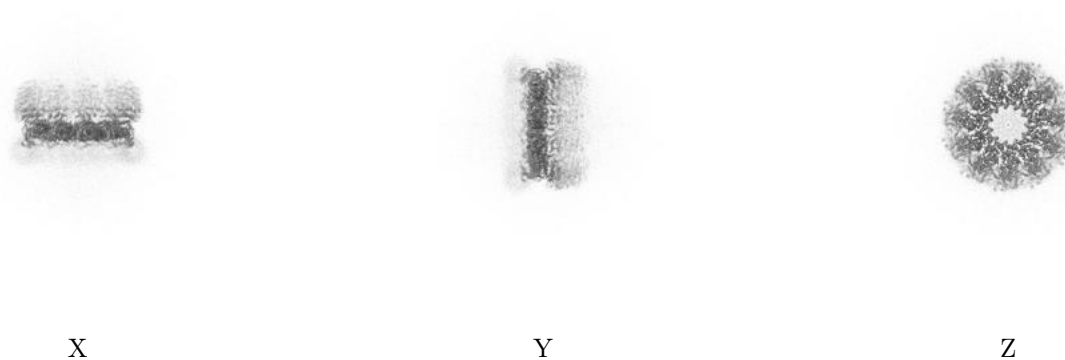
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 3.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

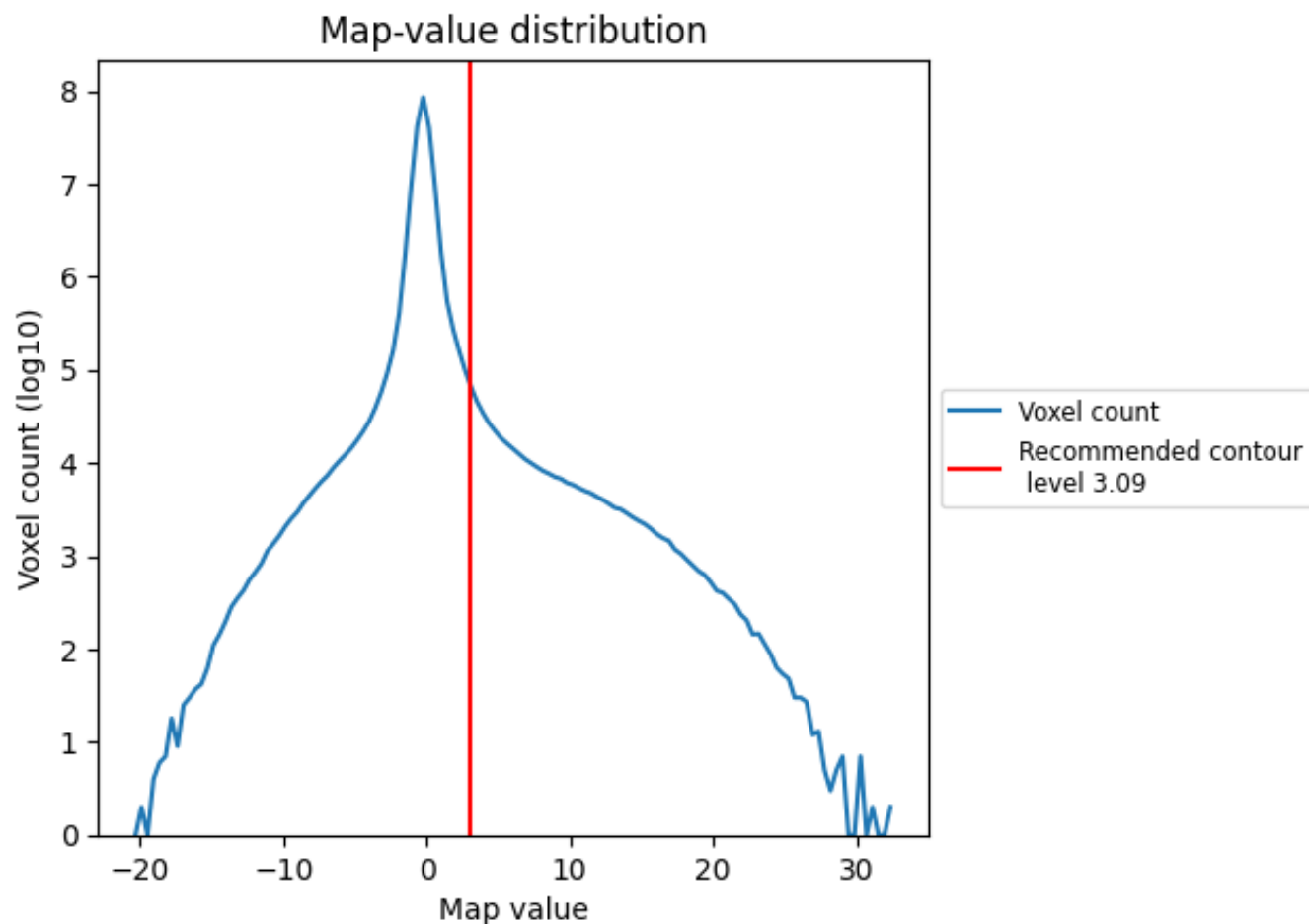
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

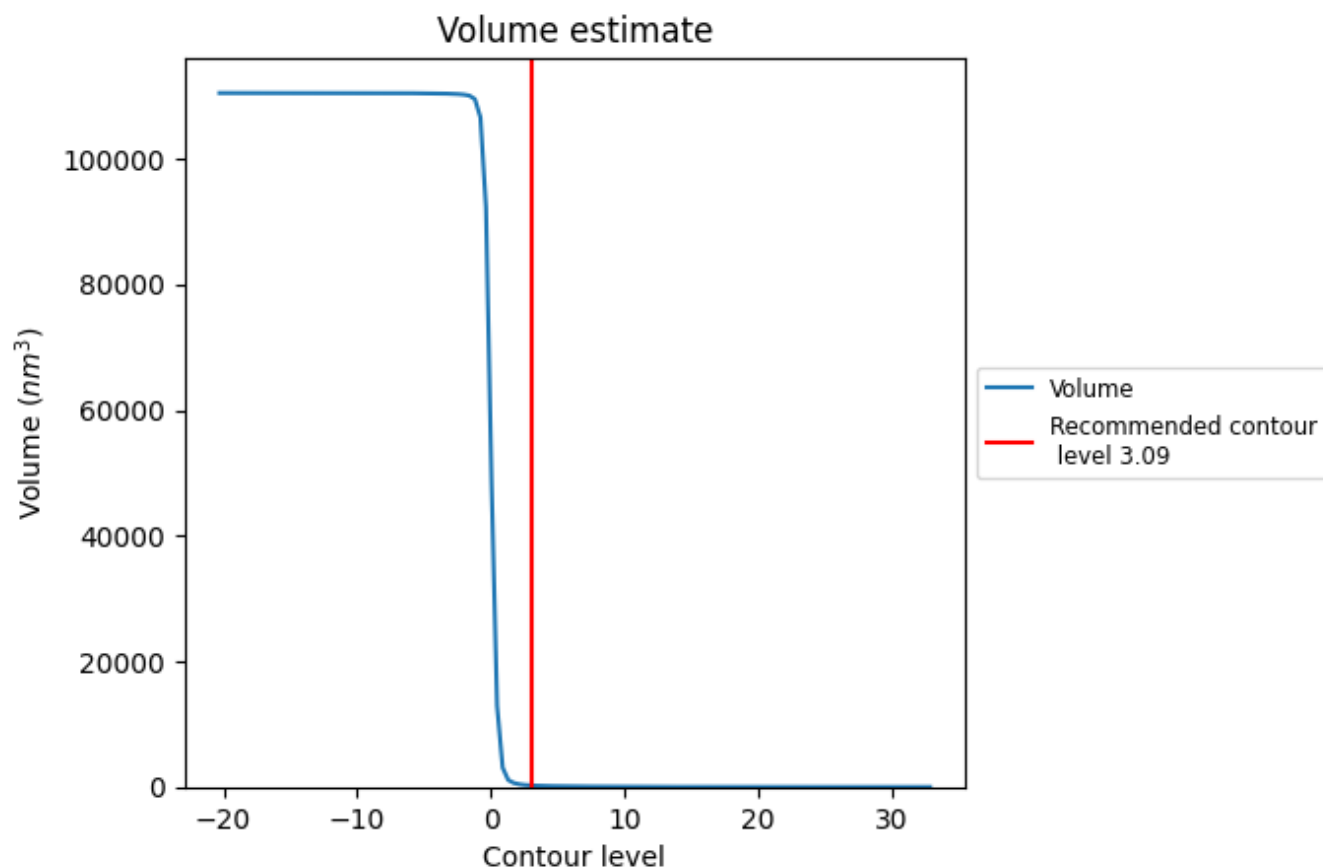
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

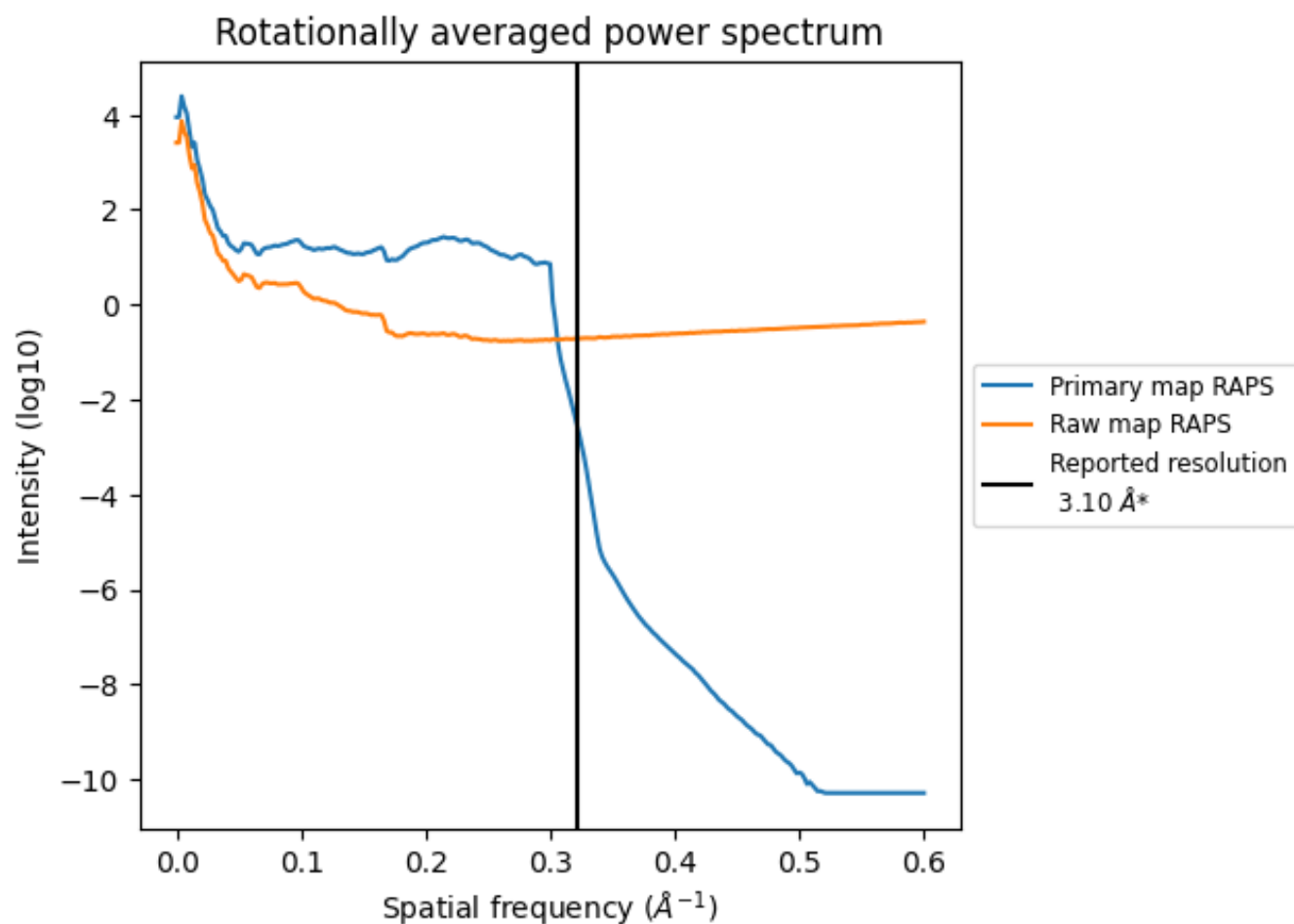
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 228 nm^3 ; this corresponds to an approximate mass of 206 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

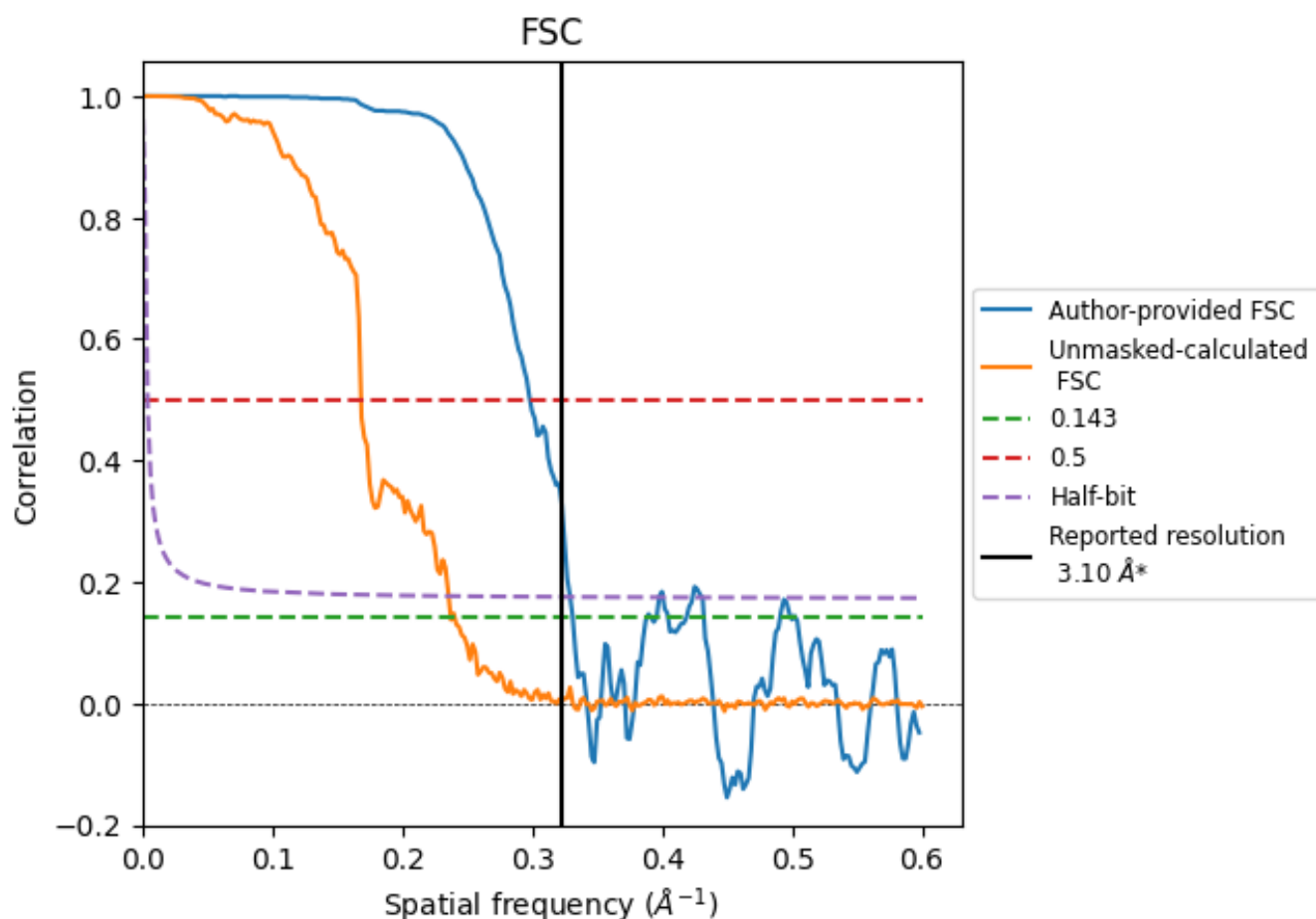


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.323 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

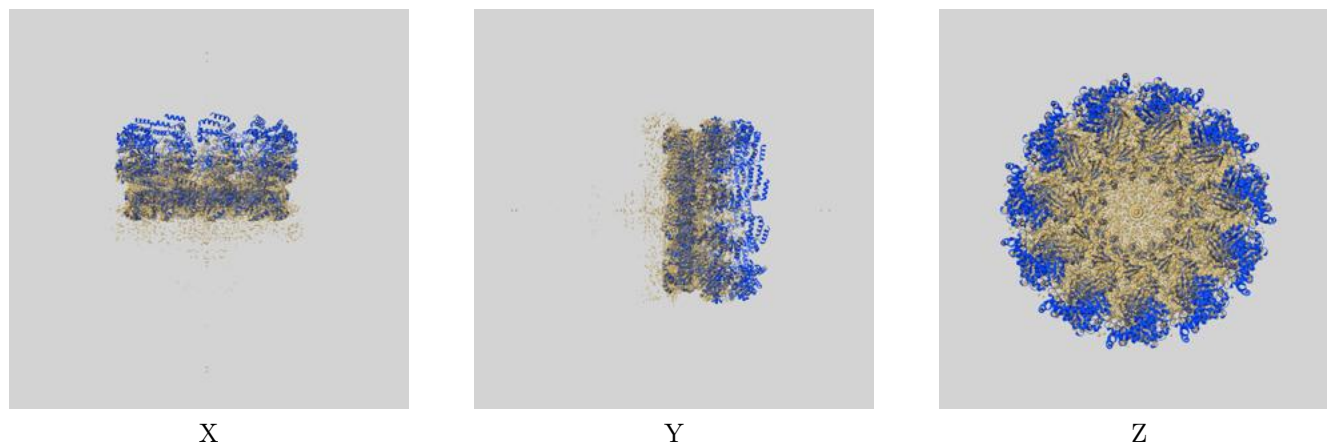
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.02	3.35	3.04
Unmasked-calculated*	4.21	5.94	4.24

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.21 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

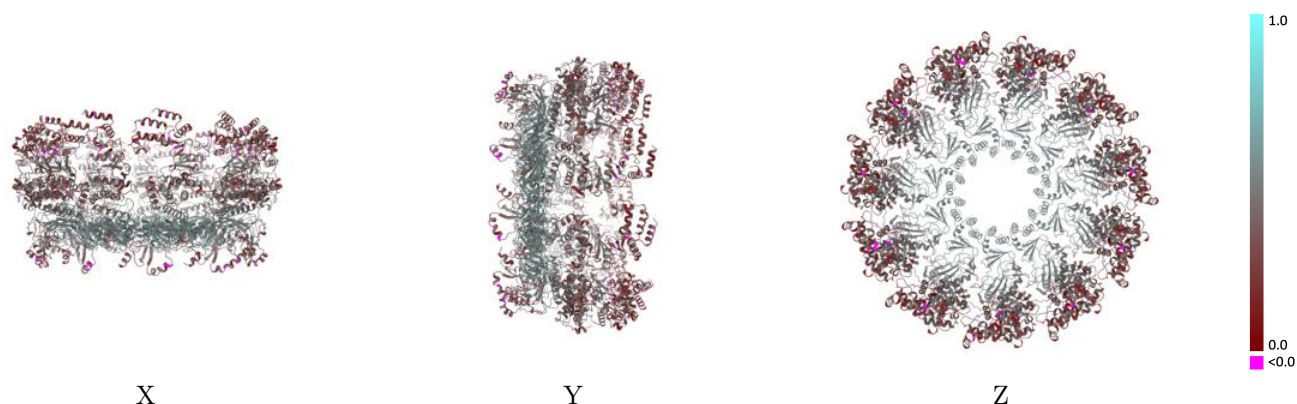
This section contains information regarding the fit between EMDB map EMD-29290 and PDB model 8FMA. Per-residue inclusion information can be found in section [3](#) on page [13](#).

9.1 Map-model overlay [i](#)



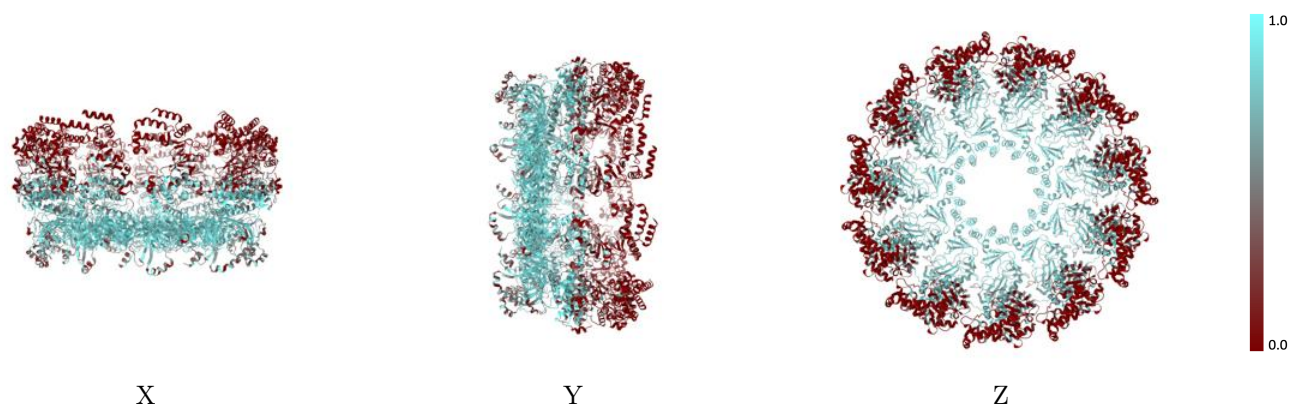
The images above show the 3D surface view of the map at the recommended contour level 3.09 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



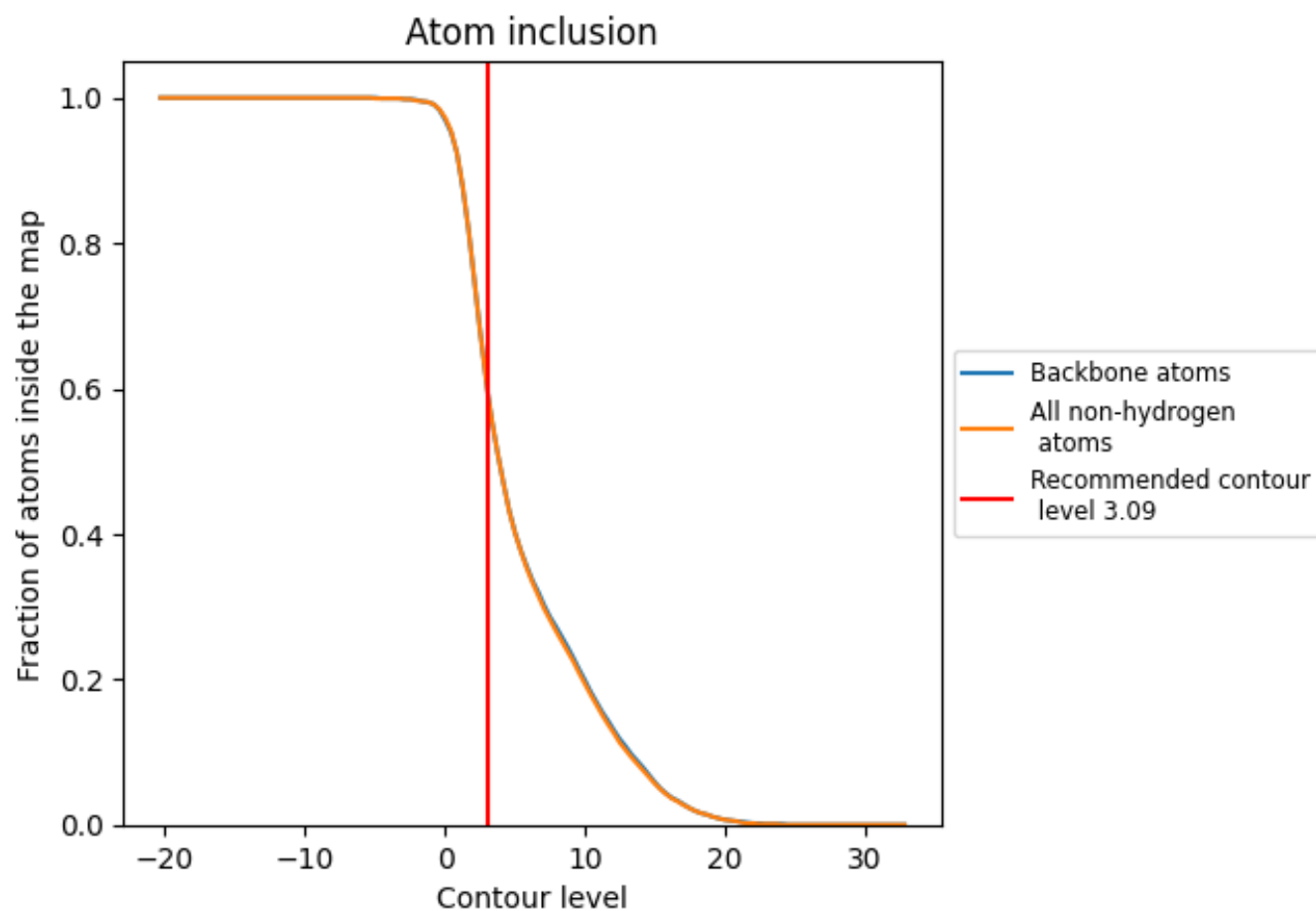
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (3.09).

9.4 Atom inclusion [i](#)



At the recommended contour level, 59% of all backbone atoms, 59% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (3.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5890	 0.4240
A	 0.7530	 0.4720
B	 0.7560	 0.4730
C	 0.7520	 0.4710
D	 0.7540	 0.4730
E	 0.7600	 0.4730
F	 0.7480	 0.4730
G	 0.7490	 0.4710
H	 0.7520	 0.4710
I	 0.7570	 0.4710
J	 0.7570	 0.4740
K	 0.7530	 0.4730
L	 0.2910	 0.3340
M	 0.2990	 0.3350
N	 0.2780	 0.3390
O	 0.2910	 0.3350
P	 0.3010	 0.3370
Q	 0.2800	 0.3400
R	 0.2880	 0.3350
S	 0.2930	 0.3320
T	 0.2870	 0.3330
U	 0.2980	 0.3360
V	 0.2990	 0.3310

