



Full wwPDB EM Validation Report ⓘ

Nov 5, 2022 – 06:15 PM EDT

PDB ID : 5W9K
EMDB ID : EMD-8786
Title : MERS S ectodomain trimer in complex with variable domain of neutralizing antibody G4
Authors : Pallesen, J.; Ward, A.B.
Deposited on : 2017-06-23
Resolution : 4.60 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.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

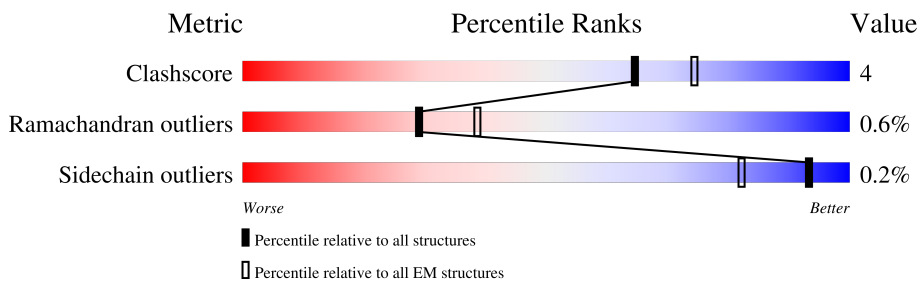
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY






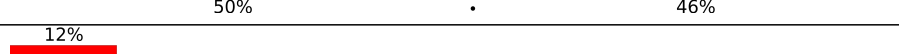


The reported resolution of this entry is 4.60 Å.

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.



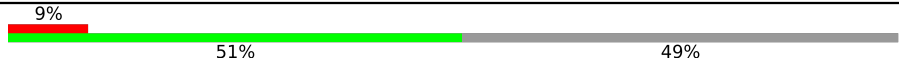


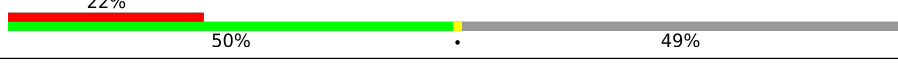
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	1329	 33% 65%
1	D	1329	 33% 65%
1	G	1329	 33% 65%
1	J	1329	 17% 51% 46%
1	K	1329	 10% 51% 46%
1	L	1329	 14% 50% 46%
2	B	233	 12% 51% 49%
2	E	233	 12% 51% 49%

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Mol	Chain	Length	Quality of chain
2	H	233	
3	C	218	
3	F	218	
3	I	218	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 32873 atoms, of which 0 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	461	3531	2233	598	683	17	0	0
1	D	462	3538	2238	599	684	17	0	0
1	G	460	3527	2231	597	682	17	0	0
1	J	724	5645	3593	924	1094	34	0	0
1	K	724	5645	3593	924	1094	34	0	0
1	L	723	5638	3589	923	1092	34	0	0

There are 258 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	506	PHE	LEU	conflict	UNP W5ZZF5
A	748	ALA	ARG	conflict	UNP W5ZZF5
A	751	GLY	ARG	conflict	UNP W5ZZF5
A	1060	PRO	VAL	conflict	UNP W5ZZF5
A	1061	PRO	LEU	conflict	UNP W5ZZF5
A	1292	GLY	-	expression tag	UNP W5ZZF5
A	1293	SER	-	expression tag	UNP W5ZZF5
A	1294	GLY	-	expression tag	UNP W5ZZF5
A	1295	TYR	-	expression tag	UNP W5ZZF5
A	1296	ILE	-	expression tag	UNP W5ZZF5
A	1297	PRO	-	expression tag	UNP W5ZZF5
A	1298	GLU	-	expression tag	UNP W5ZZF5
A	1299	ALA	-	expression tag	UNP W5ZZF5
A	1300	PRO	-	expression tag	UNP W5ZZF5
A	1301	ARG	-	expression tag	UNP W5ZZF5
A	1302	ASP	-	expression tag	UNP W5ZZF5
A	1303	GLY	-	expression tag	UNP W5ZZF5
A	1304	GLN	-	expression tag	UNP W5ZZF5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1305	ALA	-	expression tag	UNP W5ZZF5
A	1306	TYR	-	expression tag	UNP W5ZZF5
A	1307	VAL	-	expression tag	UNP W5ZZF5
A	1308	ARG	-	expression tag	UNP W5ZZF5
A	1309	LYS	-	expression tag	UNP W5ZZF5
A	1310	ASP	-	expression tag	UNP W5ZZF5
A	1311	GLY	-	expression tag	UNP W5ZZF5
A	1312	GLU	-	expression tag	UNP W5ZZF5
A	1313	TRP	-	expression tag	UNP W5ZZF5
A	1314	VAL	-	expression tag	UNP W5ZZF5
A	1315	LEU	-	expression tag	UNP W5ZZF5
A	1316	LEU	-	expression tag	UNP W5ZZF5
A	1317	SER	-	expression tag	UNP W5ZZF5
A	1318	THR	-	expression tag	UNP W5ZZF5
A	1319	PHE	-	expression tag	UNP W5ZZF5
A	1320	LEU	-	expression tag	UNP W5ZZF5
A	1321	GLY	-	expression tag	UNP W5ZZF5
A	1322	ARG	-	expression tag	UNP W5ZZF5
A	1323	SER	-	expression tag	UNP W5ZZF5
A	1324	LEU	-	expression tag	UNP W5ZZF5
A	1325	GLU	-	expression tag	UNP W5ZZF5
A	1326	VAL	-	expression tag	UNP W5ZZF5
A	1327	LEU	-	expression tag	UNP W5ZZF5
A	1328	PHE	-	expression tag	UNP W5ZZF5
A	1329	GLN	-	expression tag	UNP W5ZZF5
D	506	PHE	LEU	conflict	UNP W5ZZF5
D	748	ALA	ARG	conflict	UNP W5ZZF5
D	751	GLY	ARG	conflict	UNP W5ZZF5
D	1060	PRO	VAL	conflict	UNP W5ZZF5
D	1061	PRO	LEU	conflict	UNP W5ZZF5
D	1292	GLY	-	expression tag	UNP W5ZZF5
D	1293	SER	-	expression tag	UNP W5ZZF5
D	1294	GLY	-	expression tag	UNP W5ZZF5
D	1295	TYR	-	expression tag	UNP W5ZZF5
D	1296	ILE	-	expression tag	UNP W5ZZF5
D	1297	PRO	-	expression tag	UNP W5ZZF5
D	1298	GLU	-	expression tag	UNP W5ZZF5
D	1299	ALA	-	expression tag	UNP W5ZZF5
D	1300	PRO	-	expression tag	UNP W5ZZF5
D	1301	ARG	-	expression tag	UNP W5ZZF5
D	1302	ASP	-	expression tag	UNP W5ZZF5
D	1303	GLY	-	expression tag	UNP W5ZZF5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1304	GLN	-	expression tag	UNP W5ZZF5
D	1305	ALA	-	expression tag	UNP W5ZZF5
D	1306	TYR	-	expression tag	UNP W5ZZF5
D	1307	VAL	-	expression tag	UNP W5ZZF5
D	1308	ARG	-	expression tag	UNP W5ZZF5
D	1309	LYS	-	expression tag	UNP W5ZZF5
D	1310	ASP	-	expression tag	UNP W5ZZF5
D	1311	GLY	-	expression tag	UNP W5ZZF5
D	1312	GLU	-	expression tag	UNP W5ZZF5
D	1313	TRP	-	expression tag	UNP W5ZZF5
D	1314	VAL	-	expression tag	UNP W5ZZF5
D	1315	LEU	-	expression tag	UNP W5ZZF5
D	1316	LEU	-	expression tag	UNP W5ZZF5
D	1317	SER	-	expression tag	UNP W5ZZF5
D	1318	THR	-	expression tag	UNP W5ZZF5
D	1319	PHE	-	expression tag	UNP W5ZZF5
D	1320	LEU	-	expression tag	UNP W5ZZF5
D	1321	GLY	-	expression tag	UNP W5ZZF5
D	1322	ARG	-	expression tag	UNP W5ZZF5
D	1323	SER	-	expression tag	UNP W5ZZF5
D	1324	LEU	-	expression tag	UNP W5ZZF5
D	1325	GLU	-	expression tag	UNP W5ZZF5
D	1326	VAL	-	expression tag	UNP W5ZZF5
D	1327	LEU	-	expression tag	UNP W5ZZF5
D	1328	PHE	-	expression tag	UNP W5ZZF5
D	1329	GLN	-	expression tag	UNP W5ZZF5
G	506	PHE	LEU	conflict	UNP W5ZZF5
G	748	ALA	ARG	conflict	UNP W5ZZF5
G	751	GLY	ARG	conflict	UNP W5ZZF5
G	1060	PRO	VAL	conflict	UNP W5ZZF5
G	1061	PRO	LEU	conflict	UNP W5ZZF5
G	1292	GLY	-	expression tag	UNP W5ZZF5
G	1293	SER	-	expression tag	UNP W5ZZF5
G	1294	GLY	-	expression tag	UNP W5ZZF5
G	1295	TYR	-	expression tag	UNP W5ZZF5
G	1296	ILE	-	expression tag	UNP W5ZZF5
G	1297	PRO	-	expression tag	UNP W5ZZF5
G	1298	GLU	-	expression tag	UNP W5ZZF5
G	1299	ALA	-	expression tag	UNP W5ZZF5
G	1300	PRO	-	expression tag	UNP W5ZZF5
G	1301	ARG	-	expression tag	UNP W5ZZF5
G	1302	ASP	-	expression tag	UNP W5ZZF5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	1303	GLY	-	expression tag	UNP W5ZZF5
G	1304	GLN	-	expression tag	UNP W5ZZF5
G	1305	ALA	-	expression tag	UNP W5ZZF5
G	1306	TYR	-	expression tag	UNP W5ZZF5
G	1307	VAL	-	expression tag	UNP W5ZZF5
G	1308	ARG	-	expression tag	UNP W5ZZF5
G	1309	LYS	-	expression tag	UNP W5ZZF5
G	1310	ASP	-	expression tag	UNP W5ZZF5
G	1311	GLY	-	expression tag	UNP W5ZZF5
G	1312	GLU	-	expression tag	UNP W5ZZF5
G	1313	TRP	-	expression tag	UNP W5ZZF5
G	1314	VAL	-	expression tag	UNP W5ZZF5
G	1315	LEU	-	expression tag	UNP W5ZZF5
G	1316	LEU	-	expression tag	UNP W5ZZF5
G	1317	SER	-	expression tag	UNP W5ZZF5
G	1318	THR	-	expression tag	UNP W5ZZF5
G	1319	PHE	-	expression tag	UNP W5ZZF5
G	1320	LEU	-	expression tag	UNP W5ZZF5
G	1321	GLY	-	expression tag	UNP W5ZZF5
G	1322	ARG	-	expression tag	UNP W5ZZF5
G	1323	SER	-	expression tag	UNP W5ZZF5
G	1324	LEU	-	expression tag	UNP W5ZZF5
G	1325	GLU	-	expression tag	UNP W5ZZF5
G	1326	VAL	-	expression tag	UNP W5ZZF5
G	1327	LEU	-	expression tag	UNP W5ZZF5
G	1328	PHE	-	expression tag	UNP W5ZZF5
G	1329	GLN	-	expression tag	UNP W5ZZF5
J	506	PHE	LEU	conflict	UNP W5ZZF5
J	748	ALA	ARG	conflict	UNP W5ZZF5
J	751	GLY	ARG	conflict	UNP W5ZZF5
J	1060	PRO	VAL	conflict	UNP W5ZZF5
J	1061	PRO	LEU	conflict	UNP W5ZZF5
J	1292	GLY	-	expression tag	UNP W5ZZF5
J	1293	SER	-	expression tag	UNP W5ZZF5
J	1294	GLY	-	expression tag	UNP W5ZZF5
J	1295	TYR	-	expression tag	UNP W5ZZF5
J	1296	ILE	-	expression tag	UNP W5ZZF5
J	1297	PRO	-	expression tag	UNP W5ZZF5
J	1298	GLU	-	expression tag	UNP W5ZZF5
J	1299	ALA	-	expression tag	UNP W5ZZF5
J	1300	PRO	-	expression tag	UNP W5ZZF5
J	1301	ARG	-	expression tag	UNP W5ZZF5

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Chain	Residue	Modelled	Actual	Comment	Reference
J	1302	ASP	-	expression tag	UNP W5ZZF5
J	1303	GLY	-	expression tag	UNP W5ZZF5
J	1304	GLN	-	expression tag	UNP W5ZZF5
J	1305	ALA	-	expression tag	UNP W5ZZF5
J	1306	TYR	-	expression tag	UNP W5ZZF5
J	1307	VAL	-	expression tag	UNP W5ZZF5
J	1308	ARG	-	expression tag	UNP W5ZZF5
J	1309	LYS	-	expression tag	UNP W5ZZF5
J	1310	ASP	-	expression tag	UNP W5ZZF5
J	1311	GLY	-	expression tag	UNP W5ZZF5
J	1312	GLU	-	expression tag	UNP W5ZZF5
J	1313	TRP	-	expression tag	UNP W5ZZF5
J	1314	VAL	-	expression tag	UNP W5ZZF5
J	1315	LEU	-	expression tag	UNP W5ZZF5
J	1316	LEU	-	expression tag	UNP W5ZZF5
J	1317	SER	-	expression tag	UNP W5ZZF5
J	1318	THR	-	expression tag	UNP W5ZZF5
J	1319	PHE	-	expression tag	UNP W5ZZF5
J	1320	LEU	-	expression tag	UNP W5ZZF5
J	1321	GLY	-	expression tag	UNP W5ZZF5
J	1322	ARG	-	expression tag	UNP W5ZZF5
J	1323	SER	-	expression tag	UNP W5ZZF5
J	1324	LEU	-	expression tag	UNP W5ZZF5
J	1325	GLU	-	expression tag	UNP W5ZZF5
J	1326	VAL	-	expression tag	UNP W5ZZF5
J	1327	LEU	-	expression tag	UNP W5ZZF5
J	1328	PHE	-	expression tag	UNP W5ZZF5
J	1329	GLN	-	expression tag	UNP W5ZZF5
K	506	PHE	LEU	conflict	UNP W5ZZF5
K	748	ALA	ARG	conflict	UNP W5ZZF5
K	751	GLY	ARG	conflict	UNP W5ZZF5
K	1060	PRO	VAL	conflict	UNP W5ZZF5
K	1061	PRO	LEU	conflict	UNP W5ZZF5
K	1292	GLY	-	expression tag	UNP W5ZZF5
K	1293	SER	-	expression tag	UNP W5ZZF5
K	1294	GLY	-	expression tag	UNP W5ZZF5
K	1295	TYR	-	expression tag	UNP W5ZZF5
K	1296	ILE	-	expression tag	UNP W5ZZF5
K	1297	PRO	-	expression tag	UNP W5ZZF5
K	1298	GLU	-	expression tag	UNP W5ZZF5
K	1299	ALA	-	expression tag	UNP W5ZZF5
K	1300	PRO	-	expression tag	UNP W5ZZF5

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Chain	Residue	Modelled	Actual	Comment	Reference
K	1301	ARG	-	expression tag	UNP W5ZZF5
K	1302	ASP	-	expression tag	UNP W5ZZF5
K	1303	GLY	-	expression tag	UNP W5ZZF5
K	1304	GLN	-	expression tag	UNP W5ZZF5
K	1305	ALA	-	expression tag	UNP W5ZZF5
K	1306	TYR	-	expression tag	UNP W5ZZF5
K	1307	VAL	-	expression tag	UNP W5ZZF5
K	1308	ARG	-	expression tag	UNP W5ZZF5
K	1309	LYS	-	expression tag	UNP W5ZZF5
K	1310	ASP	-	expression tag	UNP W5ZZF5
K	1311	GLY	-	expression tag	UNP W5ZZF5
K	1312	GLU	-	expression tag	UNP W5ZZF5
K	1313	TRP	-	expression tag	UNP W5ZZF5
K	1314	VAL	-	expression tag	UNP W5ZZF5
K	1315	LEU	-	expression tag	UNP W5ZZF5
K	1316	LEU	-	expression tag	UNP W5ZZF5
K	1317	SER	-	expression tag	UNP W5ZZF5
K	1318	THR	-	expression tag	UNP W5ZZF5
K	1319	PHE	-	expression tag	UNP W5ZZF5
K	1320	LEU	-	expression tag	UNP W5ZZF5
K	1321	GLY	-	expression tag	UNP W5ZZF5
K	1322	ARG	-	expression tag	UNP W5ZZF5
K	1323	SER	-	expression tag	UNP W5ZZF5
K	1324	LEU	-	expression tag	UNP W5ZZF5
K	1325	GLU	-	expression tag	UNP W5ZZF5
K	1326	VAL	-	expression tag	UNP W5ZZF5
K	1327	LEU	-	expression tag	UNP W5ZZF5
K	1328	PHE	-	expression tag	UNP W5ZZF5
K	1329	GLN	-	expression tag	UNP W5ZZF5
L	506	PHE	LEU	conflict	UNP W5ZZF5
L	748	ALA	ARG	conflict	UNP W5ZZF5
L	751	GLY	ARG	conflict	UNP W5ZZF5
L	1060	PRO	VAL	conflict	UNP W5ZZF5
L	1061	PRO	LEU	conflict	UNP W5ZZF5
L	1292	GLY	-	expression tag	UNP W5ZZF5
L	1293	SER	-	expression tag	UNP W5ZZF5
L	1294	GLY	-	expression tag	UNP W5ZZF5
L	1295	TYR	-	expression tag	UNP W5ZZF5
L	1296	ILE	-	expression tag	UNP W5ZZF5
L	1297	PRO	-	expression tag	UNP W5ZZF5
L	1298	GLU	-	expression tag	UNP W5ZZF5
L	1299	ALA	-	expression tag	UNP W5ZZF5

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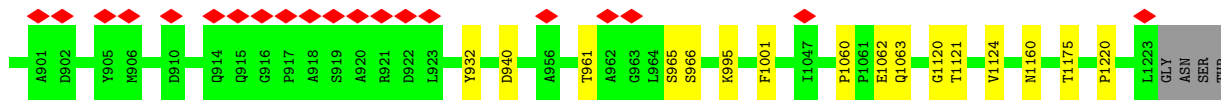
Chain	Residue	Modelled	Actual	Comment	Reference
L	1300	PRO	-	expression tag	UNP W5ZZF5
L	1301	ARG	-	expression tag	UNP W5ZZF5
L	1302	ASP	-	expression tag	UNP W5ZZF5
L	1303	GLY	-	expression tag	UNP W5ZZF5
L	1304	GLN	-	expression tag	UNP W5ZZF5
L	1305	ALA	-	expression tag	UNP W5ZZF5
L	1306	TYR	-	expression tag	UNP W5ZZF5
L	1307	VAL	-	expression tag	UNP W5ZZF5
L	1308	ARG	-	expression tag	UNP W5ZZF5
L	1309	LYS	-	expression tag	UNP W5ZZF5
L	1310	ASP	-	expression tag	UNP W5ZZF5
L	1311	GLY	-	expression tag	UNP W5ZZF5
L	1312	GLU	-	expression tag	UNP W5ZZF5
L	1313	TRP	-	expression tag	UNP W5ZZF5
L	1314	VAL	-	expression tag	UNP W5ZZF5
L	1315	LEU	-	expression tag	UNP W5ZZF5
L	1316	LEU	-	expression tag	UNP W5ZZF5
L	1317	SER	-	expression tag	UNP W5ZZF5
L	1318	THR	-	expression tag	UNP W5ZZF5
L	1319	PHE	-	expression tag	UNP W5ZZF5
L	1320	LEU	-	expression tag	UNP W5ZZF5
L	1321	GLY	-	expression tag	UNP W5ZZF5
L	1322	ARG	-	expression tag	UNP W5ZZF5
L	1323	SER	-	expression tag	UNP W5ZZF5
L	1324	LEU	-	expression tag	UNP W5ZZF5
L	1325	GLU	-	expression tag	UNP W5ZZF5
L	1326	VAL	-	expression tag	UNP W5ZZF5
L	1327	LEU	-	expression tag	UNP W5ZZF5
L	1328	PHE	-	expression tag	UNP W5ZZF5
L	1329	GLN	-	expression tag	UNP W5ZZF5

- Molecule 2 is a protein called G4 VH.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	119	Total	C	N	O	S	0	0
			948	602	156	185	5		
2	E	119	Total	C	N	O	S	0	0
			948	602	156	185	5		
2	H	119	Total	C	N	O	S	0	0
			948	602	156	185	5		

- Molecule 3 is a protein called G4 VL.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	111	Total 835	522	143	166	4	0	0
3	F	111	Total 835	522	143	166	4	0	0
3	I	111	Total 835	522	143	166	4	0	0



GLY	ILE	ASP	PHE	GLN	ASP	GLU	LEU	ASP	PHE	GLU	PHE	PHE	LYS	ASN	VAL	SER	THR	THR	ILE	PRO	ASN	PHE	GLY	SER	LEU	THR
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

ASN	TYR	THR	TYR	GLY	SER	GLY	TYR	ILE	PRO	GLU	ALA	ALA	LYS	ARG	ASN	VAL	ARG	THR	THR	THR	THR	THR	THR	THR	THR	THR	GLN
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• Molecule 1: Spike glycoprotein



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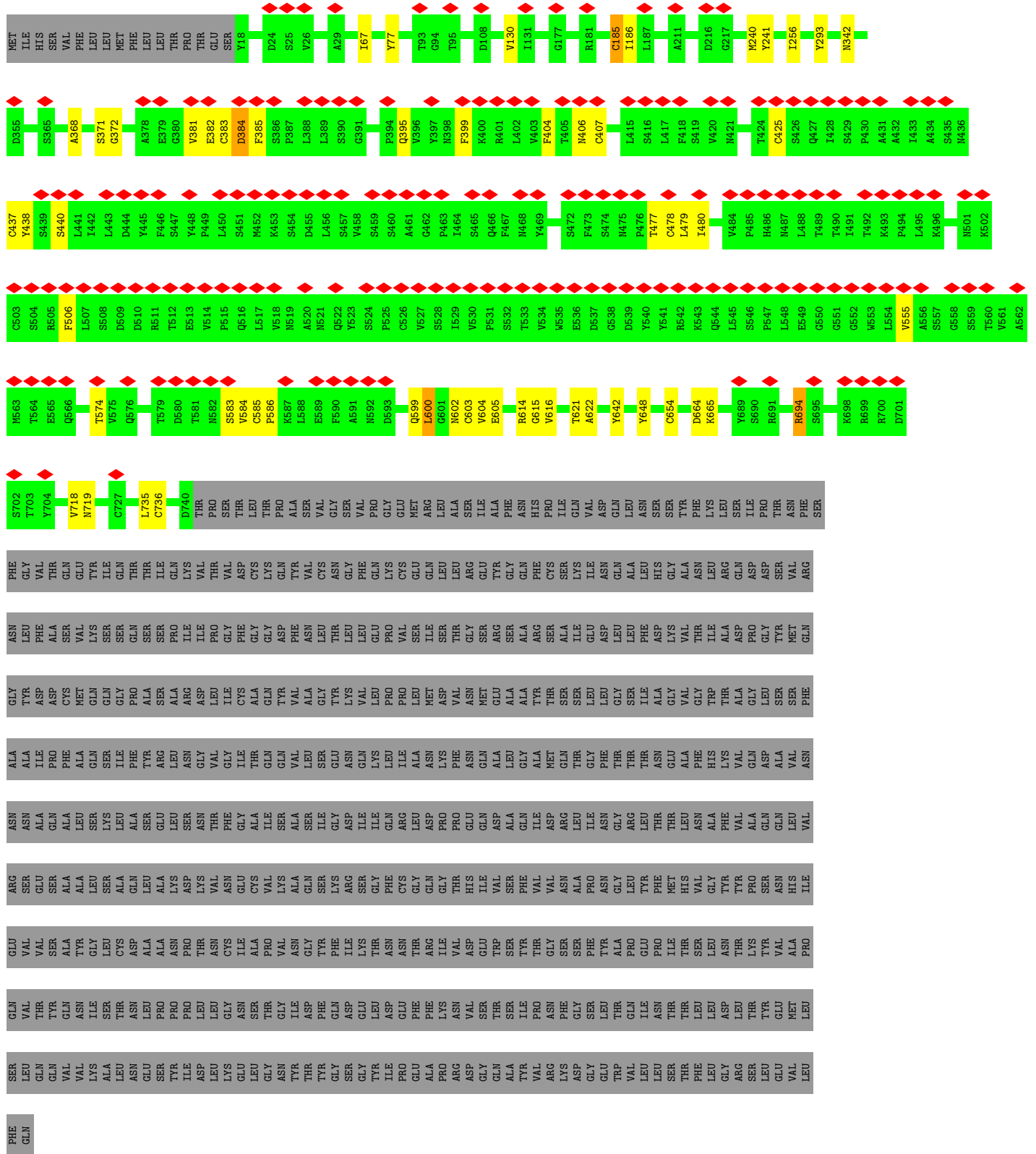
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VAL	ILE	TYR	ASP	GLU	HIS	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	SER
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SER	LEU	PHE	VAL	GLU	THR	GLY	THR	GLY	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR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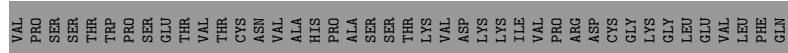
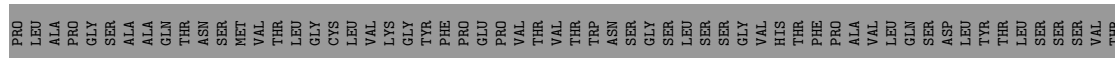
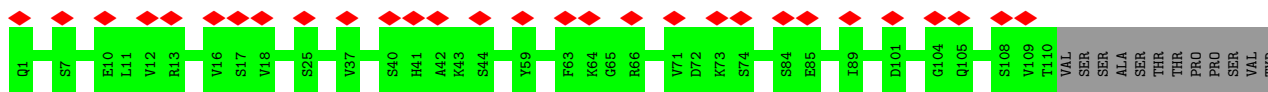


SER	LEU	PHE	VAL	GLU	THR	GLY	THR	GLY	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	SER
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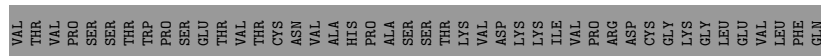
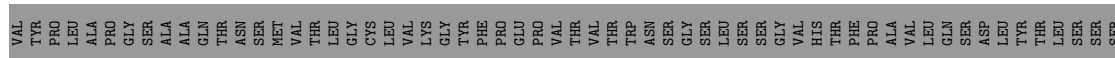
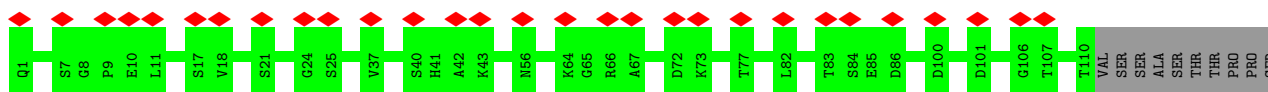


• Molecule 2: G4 VH

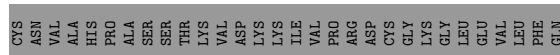
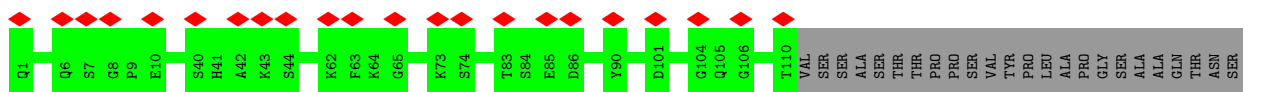




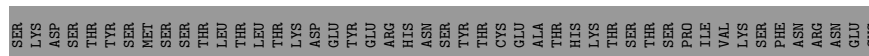
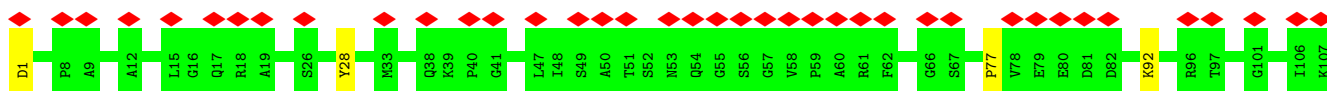
• Molecule 2: G4 VH



• Molecule 2: G4 VH

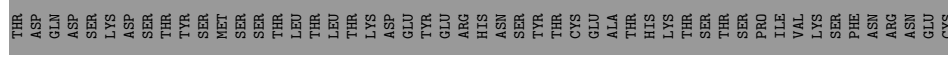
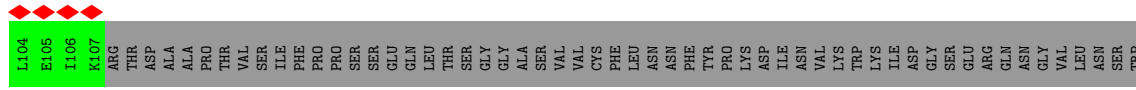
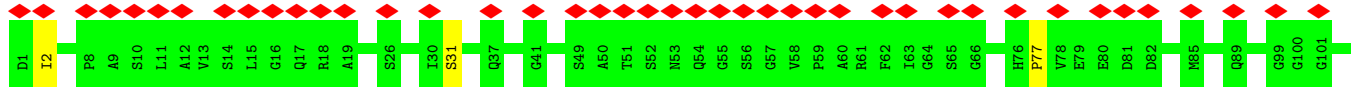


• Molecule 3: G4 VL

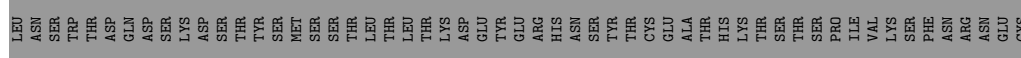
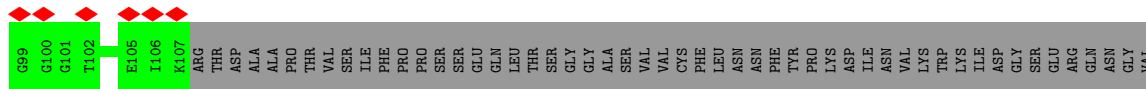
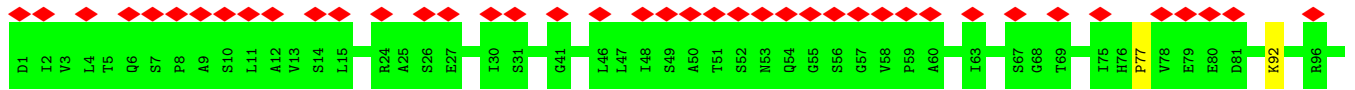


• Molecule 3: G4 VL





• Molecule 3: G4 VL



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	10544	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 ($e^-/\text{\AA}^2$)	1.89	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	29000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.066	Depositor
Minimum map value	-0.031	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.015	Depositor
Map size (Å)	310.08, 310.08, 310.08	wwPDB
Map dimensions	304, 304, 304	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.02, 1.02, 1.02	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.75	0/3603	0.82	2/4899 (0.0%)
1	D	0.76	0/3611	0.81	2/4910 (0.0%)
1	G	0.77	0/3599	0.84	2/4894 (0.0%)
1	J	0.72	0/5789	0.89	12/7881 (0.2%)
1	K	0.73	0/5789	0.89	8/7881 (0.1%)
1	L	0.71	0/5782	0.89	11/7871 (0.1%)
2	B	0.71	0/972	0.82	0/1317
2	E	0.72	0/972	0.80	0/1317
2	H	0.70	0/972	0.81	0/1317
3	C	0.72	0/852	0.77	0/1153
3	F	0.76	0/852	0.84	0/1153
3	I	0.72	0/852	0.78	0/1153
All	All	0.73	0/33645	0.86	37/45746 (0.1%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	932	TYR	CB-CG-CD1	-9.33	115.40	121.00
1	A	932	TYR	CB-CG-CD1	-8.11	116.13	121.00
1	J	642	TYR	CB-CG-CD2	-7.54	116.47	121.00
1	L	185	CYS	O-C-N	-7.32	110.99	122.70
1	L	736	CYS	N-CA-C	-7.31	91.27	111.00
1	K	184	TYR	CB-CG-CD1	-7.19	116.69	121.00
1	L	642	TYR	CB-CG-CD2	-6.72	116.97	121.00
1	L	736	CYS	CA-CB-SG	6.70	126.06	114.00
1	J	694	ARG	NE-CZ-NH1	6.59	123.59	120.30
1	L	694	ARG	NE-CZ-NH1	6.41	123.50	120.30
1	J	307	ARG	NE-CZ-NH1	-6.30	117.15	120.30
1	K	307	ARG	NE-CZ-NH2	6.24	123.42	120.30
1	J	190	ARG	NE-CZ-NH2	-6.18	117.21	120.30
1	K	323	PHE	CB-CG-CD2	-6.15	116.50	120.80
1	J	694	ARG	NE-CZ-NH2	-6.14	117.23	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	735	LEU	C-N-CA	-6.13	106.37	121.70
1	K	185	CYS	O-C-N	5.96	132.23	122.70
1	J	101	PHE	CB-CG-CD2	-5.83	116.72	120.80
1	L	694	ARG	NE-CZ-NH2	-5.67	117.46	120.30
1	L	77	TYR	CB-CG-CD2	-5.61	117.64	121.00
1	K	323	PHE	CB-CG-CD1	5.47	124.63	120.80
1	D	909	TYR	CB-CG-CD2	-5.42	117.75	121.00
1	K	378	ALA	C-N-CA	5.40	135.21	121.70
1	J	58	TYR	CB-CG-CD2	-5.37	117.78	121.00
1	K	642	TYR	CB-CG-CD1	-5.36	117.78	121.00
1	J	351	TYR	CB-CG-CD2	-5.32	117.81	121.00
1	J	438	TYR	CB-CG-CD2	-5.31	117.82	121.00
1	J	663	TYR	CB-CG-CD1	-5.26	117.84	121.00
1	G	868	ASP	CB-CG-OD2	5.23	123.01	118.30
1	D	868	ASP	CB-CG-OD2	5.23	123.00	118.30
1	L	293	TYR	CB-CG-CD2	-5.20	117.88	121.00
1	A	868	ASP	CB-CG-OD2	5.19	122.97	118.30
1	J	163	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	L	241	TYR	CB-CG-CD2	-5.10	117.94	121.00
1	L	585	CYS	C-N-CD	5.06	139.03	128.40
1	K	738	LEU	C-N-CD	5.06	139.02	128.40
1	J	585	CYS	C-N-CD	5.03	138.96	128.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3531	0	3454	15	0
1	D	3538	0	3463	20	0
1	G	3527	0	3453	14	0
1	J	5645	0	5413	74	0
1	K	5645	0	5413	52	0
1	L	5638	0	5408	91	0
2	B	948	0	904	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	948	0	904	0	0
2	H	948	0	904	0	0
3	C	835	0	816	3	0
3	F	835	0	816	1	0
3	I	835	0	816	1	0
All	All	32873	0	31764	270	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 4.

All (270) 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:L:371:SER:CB	1:L:604:VAL:HG12	1.25	1.63
1:J:371:SER:CB	1:J:604:VAL:HG12	1.17	1.59
1:L:506:PHE:CE2	1:L:555:VAL:CG2	1.77	1.54
1:L:506:PHE:CE2	1:L:555:VAL:HG21	0.85	1.37
1:L:506:PHE:HE2	1:L:555:VAL:CG2	1.20	1.36
1:J:371:SER:HB3	1:J:604:VAL:CG1	1.52	1.34
1:J:371:SER:CB	1:J:604:VAL:CG1	2.01	1.33
1:D:905:TYR:OH	1:D:936:PRO:CA	1.78	1.31
1:L:371:SER:HB3	1:L:604:VAL:CG1	1.60	1.31
1:L:384:ASP:O	1:L:404:PHE:CZ	1.83	1.31
1:L:506:PHE:CD2	1:L:555:VAL:HG21	1.69	1.27
1:L:371:SER:CB	1:L:604:VAL:CG1	2.10	1.26
1:L:506:PHE:CD2	1:L:555:VAL:CG2	2.22	1.21
1:D:905:TYR:OH	1:D:936:PRO:HA	1.47	1.13
1:J:371:SER:OG	1:J:604:VAL:CG1	1.96	1.13
1:K:477:THR:OG1	1:K:573:ILE:O	1.65	1.10
1:J:185:CYS:HB2	1:J:237:CYS:HA	1.34	1.09
1:L:371:SER:OG	1:L:604:VAL:CG1	1.99	1.09
1:K:506:PHE:CE1	1:K:513:GLU:OE2	2.06	1.08
1:D:905:TYR:OH	1:D:936:PRO:N	1.87	1.07
1:J:371:SER:OG	1:J:604:VAL:HG12	1.52	1.02
1:L:371:SER:OG	1:L:604:VAL:HG12	1.55	1.02
1:J:371:SER:CA	1:J:604:VAL:HG12	1.89	1.01
1:A:868:ASP:OD2	1:A:995:LYS:NZ	1.94	0.99
1:J:185:CYS:SG	1:J:186:ILE:N	2.29	0.99
1:J:187:LEU:HD21	1:J:228:PHE:HZ	1.25	0.99
1:L:371:SER:HB3	1:L:604:VAL:HG12	0.99	0.97
1:L:384:ASP:CB	1:L:404:PHE:HE2	1.75	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:384:ASP:O	1:L:404:PHE:CE2	2.18	0.97
1:L:621:THR:O	1:L:648:TYR:HD2	1.45	0.96
1:J:185:CYS:SG	1:J:235:ARG:O	2.23	0.96
1:L:372:GLY:O	1:L:604:VAL:HB	1.66	0.95
1:J:371:SER:HB3	1:J:604:VAL:HG12	0.98	0.95
1:D:905:TYR:CZ	1:D:936:PRO:HB3	2.03	0.92
1:L:506:PHE:CD2	1:L:555:VAL:HG23	2.03	0.92
1:G:912:CYS:HG	1:G:925:CYS:HG	1.09	0.91
1:J:185:CYS:CB	1:J:237:CYS:HA	2.00	0.90
1:L:371:SER:CA	1:L:604:VAL:HG12	2.02	0.90
1:J:185:CYS:HB2	1:J:237:CYS:CA	2.02	0.89
1:J:372:GLY:O	1:J:604:VAL:HB	1.74	0.87
1:J:184:TYR:CE1	1:J:286:VAL:HG11	2.11	0.86
1:J:185:CYS:HG	1:J:237:CYS:HG	1.10	0.84
1:L:425:CYS:SG	1:L:478:CYS:CB	2.66	0.84
1:K:477:THR:HB	1:K:574:THR:HB	1.60	0.83
1:L:506:PHE:HE2	1:L:555:VAL:HG22	1.40	0.83
1:L:621:THR:O	1:L:648:TYR:CD2	2.30	0.83
1:J:371:SER:HB3	1:J:604:VAL:HG13	1.57	0.83
1:K:471:GLN:HG2	1:K:477:THR:HG21	1.61	0.82
1:L:425:CYS:CB	1:L:478:CYS:SG	2.68	0.81
1:L:425:CYS:SG	1:L:478:CYS:HB3	2.19	0.81
1:L:425:CYS:HB3	1:L:478:CYS:SG	2.21	0.81
1:L:384:ASP:O	1:L:404:PHE:HZ	1.58	0.80
1:K:471:GLN:CG	1:K:477:THR:HG21	2.13	0.79
1:K:477:THR:HB	1:K:574:THR:CB	2.11	0.79
1:J:184:TYR:OH	1:J:288:ASP:OD2	2.01	0.79
1:J:184:TYR:CG	1:J:286:VAL:HG21	2.18	0.78
1:L:372:GLY:O	1:L:604:VAL:CB	2.31	0.78
1:D:905:TYR:HH	1:D:936:PRO:HA	1.49	0.77
1:L:477:THR:HB	1:L:574:THR:HG22	1.65	0.77
1:K:506:PHE:CZ	1:K:513:GLU:OE2	2.37	0.76
1:L:602:ASN:O	1:L:616:VAL:HB	1.84	0.76
1:J:187:LEU:HD21	1:J:228:PHE:CZ	2.17	0.75
1:L:384:ASP:CB	1:L:404:PHE:CE2	2.65	0.75
1:L:425:CYS:CB	1:L:478:CYS:HG	2.00	0.74
1:J:371:SER:C	1:J:604:VAL:HG12	2.08	0.74
1:K:477:THR:CB	1:K:573:ILE:O	2.35	0.74
1:K:369:LYS:O	1:K:369:LYS:HG2	1.88	0.74
1:D:905:TYR:OH	1:D:936:PRO:CB	2.36	0.72
1:J:372:GLY:O	1:J:604:VAL:CB	2.37	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:381:VAL:HG11	1:L:407:CYS:HA	1.70	0.72
1:L:621:THR:HG22	1:L:622:ALA:N	2.03	0.72
1:D:905:TYR:CZ	1:D:936:PRO:CB	2.73	0.72
1:J:602:ASN:O	1:J:616:VAL:HB	1.90	0.71
1:L:384:ASP:HB2	1:L:404:PHE:HE2	1.54	0.70
1:L:371:SER:HB3	1:L:604:VAL:HG13	1.71	0.70
1:L:477:THR:CG2	1:L:574:THR:HG22	2.21	0.69
1:J:185:CYS:CB	1:J:237:CYS:SG	2.81	0.69
1:J:185:CYS:CB	1:J:237:CYS:HG	2.05	0.69
1:L:506:PHE:HD2	1:L:555:VAL:HG23	1.56	0.68
1:L:406:ASN:HA	1:L:583:SER:HB2	1.76	0.67
1:J:185:CYS:HB2	1:J:237:CYS:CB	2.26	0.66
1:K:506:PHE:CD1	1:K:513:GLU:OE2	2.48	0.66
1:J:371:SER:OG	1:J:604:VAL:HG11	1.92	0.66
1:L:477:THR:CB	1:L:574:THR:HG22	2.26	0.65
1:K:477:THR:HG1	1:K:478:CYS:H	1.45	0.65
1:K:477:THR:HG1	1:K:573:ILE:C	1.92	0.65
1:K:471:GLN:CD	1:K:477:THR:HG21	2.17	0.64
1:L:371:SER:C	1:L:604:VAL:HG12	2.17	0.64
1:J:184:TYR:CD1	1:J:286:VAL:HG21	2.31	0.64
1:K:621:THR:O	1:K:648:TYR:HD2	1.80	0.64
1:K:621:THR:HG22	1:K:622:ALA:N	2.11	0.64
1:L:425:CYS:HG	1:L:478:CYS:CB	2.06	0.64
1:D:905:TYR:CZ	1:D:936:PRO:CA	2.80	0.64
1:J:184:TYR:HB3	1:J:239:PHE:HB2	1.80	0.63
1:K:383:CYS:SG	1:K:409:TYR:HB3	2.38	0.63
1:J:186:ILE:HG22	1:J:187:LEU:N	2.13	0.63
1:L:384:ASP:CA	1:L:404:PHE:HE2	2.12	0.62
1:L:382:GLU:HA	1:L:382:GLU:OE1	1.99	0.62
1:A:897:LYS:HE2	1:A:897:LYS:HA	1.81	0.62
1:J:184:TYR:HD2	1:J:239:PHE:HD1	1.45	0.62
1:J:185:CYS:SG	1:J:237:CYS:HA	2.39	0.61
1:L:384:ASP:HB3	1:L:404:PHE:HE2	1.65	0.61
1:J:184:TYR:CZ	1:J:286:VAL:CG1	2.83	0.61
1:L:384:ASP:HB3	1:L:404:PHE:CE2	2.34	0.61
1:L:477:THR:HG22	1:L:574:THR:HB	1.83	0.61
1:K:738:LEU:HD12	1:K:739:PRO:HD3	1.82	0.60
1:J:184:TYR:CZ	1:J:286:VAL:HG13	2.37	0.60
1:L:621:THR:CG2	1:L:622:ALA:N	2.64	0.59
1:L:371:SER:OG	1:L:604:VAL:HG11	1.97	0.59
1:J:616:VAL:HG13	1:J:616:VAL:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:184:TYR:CD2	1:J:239:PHE:HD1	2.21	0.58
1:L:599:GLN:C	1:L:600:LEU:HG	2.24	0.58
1:L:602:ASN:O	1:L:616:VAL:CB	2.52	0.58
1:J:184:TYR:CD2	1:J:286:VAL:CG2	2.88	0.57
1:L:616:VAL:O	1:L:616:VAL:HG13	2.03	0.57
1:J:371:SER:C	1:J:604:VAL:CG1	2.73	0.56
1:J:599:GLN:C	1:J:600:LEU:HG	2.25	0.56
1:J:184:TYR:CE2	1:J:286:VAL:HG22	2.41	0.56
1:J:187:LEU:CD2	1:J:228:PHE:HZ	2.09	0.56
1:L:506:PHE:CE2	1:L:555:VAL:HG22	2.20	0.56
1:J:372:GLY:O	1:J:604:VAL:CG2	2.54	0.56
1:J:602:ASN:O	1:J:616:VAL:CB	2.54	0.56
1:K:478:CYS:O	1:K:479:LEU:HG	2.06	0.55
1:D:1175:THR:HG23	1:D:1175:THR:O	2.07	0.55
1:D:905:TYR:CZ	1:D:936:PRO:HA	2.40	0.55
1:K:621:THR:CG2	1:K:622:ALA:N	2.70	0.54
1:L:384:ASP:CA	1:L:404:PHE:CE2	2.91	0.54
1:J:184:TYR:CE1	1:J:286:VAL:CG1	2.86	0.53
1:G:1175:THR:HG23	1:G:1175:THR:O	2.07	0.53
1:A:1175:THR:HG23	1:A:1175:THR:O	2.09	0.53
1:J:184:TYR:HD2	1:J:239:PHE:CD1	2.24	0.53
1:L:384:ASP:HB2	1:L:404:PHE:CE2	2.40	0.53
1:L:67:ILE:HG13	1:L:67:ILE:O	2.08	0.53
1:K:172:LEU:O	1:K:172:LEU:HG	2.08	0.53
1:G:875:GLU:OE1	1:G:875:GLU:N	2.35	0.53
1:A:875:GLU:OE1	1:A:875:GLU:N	2.36	0.53
1:K:477:THR:HA	1:K:574:THR:HA	1.91	0.53
1:L:381:VAL:CG1	1:L:407:CYS:HA	2.39	0.52
1:L:621:THR:CG2	1:L:622:ALA:H	2.22	0.52
1:J:185:CYS:SG	1:J:237:CYS:CA	2.98	0.52
1:J:184:TYR:CD2	1:J:286:VAL:HG22	2.45	0.51
1:K:738:LEU:HD12	1:K:739:PRO:CD	2.41	0.51
1:A:875:GLU:O	1:A:875:GLU:HG2	2.09	0.51
1:L:399:PHE:O	1:L:399:PHE:CD1	2.64	0.51
1:G:826:GLN:OE1	1:G:826:GLN:N	2.27	0.50
1:K:477:THR:HG1	1:K:478:CYS:N	2.09	0.50
1:L:240:MET:SD	1:L:240:MET:C	2.89	0.50
1:L:256:ILE:HG23	1:L:256:ILE:O	2.11	0.49
1:J:184:TYR:CD2	1:J:286:VAL:HG21	2.47	0.49
1:L:477:THR:CG2	1:L:574:THR:CG2	2.90	0.49
1:G:912:CYS:SG	1:K:655:VAL:HG12	2.52	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:477:THR:OG1	1:K:478:CYS:N	2.44	0.49
1:L:372:GLY:O	1:L:604:VAL:CG2	2.60	0.49
1:J:166:ASN:O	1:J:186:ILE:HG23	2.11	0.49
1:L:371:SER:HG	1:L:604:VAL:CG1	2.23	0.49
1:L:384:ASP:C	1:L:404:PHE:CE2	2.85	0.48
1:A:1124:VAL:HG13	1:A:1124:VAL:O	2.13	0.48
1:G:897:LYS:HE2	1:G:897:LYS:HA	1.95	0.48
1:J:186:ILE:CG2	1:J:187:LEU:N	2.75	0.48
1:K:471:GLN:HG2	1:K:477:THR:CG2	2.38	0.48
1:L:477:THR:HG22	1:L:574:THR:CB	2.42	0.48
1:J:186:ILE:C	1:J:187:LEU:HD12	2.34	0.48
1:K:477:THR:CA	1:K:573:ILE:O	2.61	0.48
1:G:875:GLU:O	1:G:875:GLU:HG2	2.13	0.48
1:L:371:SER:CB	1:L:604:VAL:HG13	2.29	0.48
1:L:437:CYS:SG	1:L:584:VAL:O	2.72	0.48
1:L:395:GLN:HA	1:L:395:GLN:OE1	2.14	0.47
1:J:694:ARG:HA	1:J:694:ARG:NE	2.29	0.47
1:J:605:GLU:HA	1:J:614:ARG:HA	1.96	0.47
1:L:382:GLU:O	1:L:407:CYS:HB2	2.15	0.47
1:L:605:GLU:HA	1:L:614:ARG:HA	1.96	0.47
1:L:694:ARG:NE	1:L:694:ARG:HA	2.29	0.47
1:L:694:ARG:HA	1:L:694:ARG:HE	1.80	0.47
1:J:372:GLY:O	1:J:604:VAL:HG21	2.14	0.47
1:K:395:GLN:OE1	1:K:395:GLN:HA	2.14	0.47
1:L:342:ASN:C	1:L:342:ASN:OD1	2.53	0.47
1:J:142:LYS:NZ	1:J:142:LYS:HB3	2.29	0.47
1:L:438:TYR:O	1:L:584:VAL:HB	2.14	0.47
1:K:478:CYS:N	1:K:573:ILE:O	2.40	0.46
1:K:439:SER:HA	1:K:582:ASN:HA	1.98	0.46
1:J:130:VAL:O	1:J:130:VAL:HG12	2.16	0.46
1:J:185:CYS:CB	1:J:237:CYS:CA	2.79	0.46
1:K:731:LEU:HB2	1:K:735:LEU:HB3	1.97	0.46
1:A:806:CYS:SG	1:A:806:CYS:O	2.74	0.46
1:J:166:ASN:O	1:J:186:ILE:HG12	2.15	0.46
1:K:606:TYR:CD1	1:K:606:TYR:C	2.88	0.46
1:G:777:TYR:N	1:G:777:TYR:CD2	2.84	0.45
1:D:875:GLU:HB3	1:D:886:ALA:HB3	1.97	0.45
1:K:477:THR:HB	1:K:574:THR:CA	2.45	0.45
1:J:371:SER:CB	1:J:604:VAL:HG13	2.21	0.45
1:K:523:TYR:N	1:K:523:TYR:CD1	2.84	0.45
1:J:665:LYS:HD3	1:J:665:LYS:C	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:621:THR:O	1:K:648:TYR:CD2	2.65	0.45
1:L:615:GLY:HA2	1:L:654:CYS:SG	2.57	0.45
1:D:806:CYS:O	1:D:806:CYS:SG	2.74	0.45
1:J:184:TYR:CD2	1:J:239:PHE:CD1	3.03	0.45
1:K:130:VAL:HG12	1:K:130:VAL:O	2.17	0.45
3:C:1:ASP:C	3:C:1:ASP:OD1	2.56	0.45
1:D:905:TYR:OH	1:D:935:LEU:C	2.54	0.45
1:D:905:TYR:CE2	1:D:936:PRO:HB3	2.51	0.44
1:D:1036:LEU:HA	1:D:1036:LEU:HD23	1.74	0.44
1:K:378:ALA:HB1	1:K:379:GLU:HA	1.99	0.44
1:L:240:MET:SD	1:L:240:MET:O	2.75	0.44
1:G:1001:PHE:CD2	1:G:1001:PHE:C	2.91	0.44
1:J:615:GLY:HA2	1:J:654:CYS:SG	2.58	0.44
1:L:621:THR:HG22	1:L:622:ALA:H	1.76	0.44
1:L:440:SER:HA	1:L:584:VAL:HG21	2.00	0.44
1:A:1062:GLU:N	1:A:1062:GLU:OE1	2.51	0.44
1:A:1160:ASN:OD1	1:A:1160:ASN:C	2.56	0.44
3:F:2:ILE:N	3:F:2:ILE:HD12	2.32	0.44
1:K:172:LEU:N	1:K:172:LEU:HD23	2.33	0.44
1:J:440:SER:HA	1:J:584:VAL:HG21	2.00	0.43
1:K:590:PHE:O	1:K:590:PHE:CG	2.72	0.43
1:J:438:TYR:O	1:J:584:VAL:HB	2.18	0.43
1:G:932:TYR:HD1	1:G:932:TYR:HA	1.45	0.43
1:A:826:GLN:OE1	1:A:826:GLN:N	2.31	0.43
1:A:965:SER:OG	1:A:966:SER:N	2.52	0.42
1:G:806:CYS:O	1:G:806:CYS:SG	2.77	0.42
1:L:479:LEU:C	1:L:480:ILE:HG13	2.40	0.42
1:L:664:ASP:OD1	1:L:664:ASP:C	2.57	0.42
1:G:1060:PRO:HA	1:G:1063:GLN:HB3	2.00	0.42
1:K:732:GLY:O	1:K:735:LEU:HB2	2.19	0.42
1:K:596:ILE:HD12	1:K:596:ILE:C	2.39	0.42
1:J:49:ASP:C	1:J:49:ASP:OD1	2.58	0.42
1:A:1001:PHE:CD2	1:A:1001:PHE:C	2.93	0.42
1:A:1060:PRO:HA	1:A:1063:GLN:HB3	2.00	0.42
1:A:1120:GLY:O	1:A:1121:THR:OG1	2.31	0.42
1:G:1124:VAL:O	1:G:1124:VAL:HG13	2.19	0.42
1:J:371:SER:C	1:J:604:VAL:HB	2.40	0.42
1:K:633:ASP:C	1:K:633:ASP:OD1	2.58	0.42
3:C:28:TYR:N	3:C:28:TYR:CD1	2.86	0.42
1:G:965:SER:OG	1:G:966:SER:N	2.52	0.42
1:A:872:THR:O	1:A:872:THR:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1173:ILE:HG13	1:D:1185:SER:OG	2.20	0.42
1:J:395:GLN:HA	1:J:395:GLN:OE1	2.20	0.42
1:J:371:SER:O	1:J:604:VAL:HB	2.19	0.42
1:L:616:VAL:O	1:L:616:VAL:CG1	2.68	0.42
1:L:185:CYS:SG	1:L:186:ILE:N	2.93	0.41
1:L:506:PHE:CE2	1:L:555:VAL:CB	2.88	0.41
1:K:172:LEU:HD23	1:K:172:LEU:H	1.85	0.41
1:K:326:ASP:C	1:K:326:ASP:OD1	2.56	0.41
1:K:409:TYR:CD1	1:K:409:TYR:C	2.94	0.41
1:D:777:TYR:N	1:D:777:TYR:CD2	2.88	0.41
1:L:371:SER:C	1:L:604:VAL:CG1	2.87	0.41
3:C:92:LYS:HD3	3:C:92:LYS:C	2.41	0.41
1:J:67:ILE:O	1:J:67:ILE:HG23	2.19	0.41
1:L:665:LYS:HD3	1:L:665:LYS:C	2.40	0.41
1:J:689:TYR:CD1	1:J:689:TYR:C	2.94	0.41
1:D:809:TYR:CD2	1:D:809:TYR:C	2.94	0.41
3:I:92:LYS:HD3	3:I:92:LYS:C	2.41	0.41
1:J:184:TYR:CE2	1:J:286:VAL:CG2	3.04	0.41
1:J:603:CYS:O	1:J:603:CYS:SG	2.79	0.41
1:K:130:VAL:O	1:K:130:VAL:CG1	2.69	0.41
1:K:209:THR:O	1:K:209:THR:HG22	2.20	0.41
1:K:621:THR:CG2	1:K:622:ALA:H	2.33	0.41
1:L:130:VAL:O	1:L:130:VAL:HG12	2.20	0.41
1:L:477:THR:O	1:L:478:CYS:SG	2.79	0.41
1:K:677:VAL:HG22	1:K:677:VAL:O	2.20	0.41
1:L:477:THR:C	1:L:478:CYS:SG	3.00	0.41
1:L:719:ASN:OD1	1:L:719:ASN:C	2.59	0.41
1:D:1124:VAL:O	1:D:1124:VAL:HG13	2.21	0.41
1:L:372:GLY:C	1:L:604:VAL:HB	2.39	0.40
1:D:775:SER:OG	1:D:776:SER:N	2.53	0.40
1:L:603:CYS:SG	1:L:603:CYS:O	2.79	0.40
1:J:130:VAL:O	1:J:130:VAL:CG1	2.68	0.40
1:K:427:GLN:HB2	1:K:476:PRO:HB3	2.04	0.40
1:L:604:VAL:O	1:L:604:VAL:HG23	2.21	0.40
1:J:372:GLY:C	1:J:604:VAL:HB	2.39	0.40
1:K:478:CYS:HB3	1:K:479:LEU:H	1.73	0.40
1:K:670:HIS:N	1:K:670:HIS:CD2	2.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	457/1329 (34%)	439 (96%)	15 (3%)	3 (1%)	22	62
1	D	458/1329 (34%)	441 (96%)	12 (3%)	5 (1%)	14	52
1	G	456/1329 (34%)	438 (96%)	15 (3%)	3 (1%)	22	62
1	J	722/1329 (54%)	685 (95%)	34 (5%)	3 (0%)	34	72
1	K	722/1329 (54%)	688 (95%)	30 (4%)	4 (1%)	25	65
1	L	721/1329 (54%)	685 (95%)	32 (4%)	4 (1%)	25	65
2	B	117/233 (50%)	115 (98%)	2 (2%)	0	100	100
2	E	117/233 (50%)	115 (98%)	2 (2%)	0	100	100
2	H	117/233 (50%)	115 (98%)	2 (2%)	0	100	100
3	C	109/218 (50%)	104 (95%)	4 (4%)	1 (1%)	17	56
3	F	109/218 (50%)	105 (96%)	2 (2%)	2 (2%)	8	42
3	I	109/218 (50%)	104 (95%)	4 (4%)	1 (1%)	17	56
All	All	4214/9327 (45%)	4034 (96%)	154 (4%)	26 (1%)	29	65

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	940	ASP
1	A	961	THR
1	D	961	THR
1	D	965	SER
1	G	940	ASP
1	J	718	VAL
1	K	718	VAL
1	L	718	VAL
3	C	77	PRO
1	D	905	TYR
1	G	961	THR
3	I	77	PRO

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Mol	Chain	Res	Type
1	A	1220	PRO
1	D	1220	PRO
3	F	31	SER
3	F	77	PRO
1	G	1220	PRO
1	K	378	ALA
1	D	956	ALA
1	J	368	ALA
1	K	629	ARG
1	L	368	ALA
1	K	44	TRP
1	L	385	PHE
1	J	586	PRO
1	L	586	PRO

5.3.2 Protein sidechains [i](#)

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	386/1148 (34%)	386 (100%)	0	100	100
1	D	387/1148 (34%)	387 (100%)	0	100	100
1	G	386/1148 (34%)	386 (100%)	0	100	100
1	J	633/1148 (55%)	631 (100%)	2 (0%)	92	95
1	K	633/1148 (55%)	631 (100%)	2 (0%)	92	95
1	L	632/1148 (55%)	629 (100%)	3 (0%)	88	93
2	B	102/202 (50%)	102 (100%)	0	100	100
2	E	102/202 (50%)	102 (100%)	0	100	100
2	H	102/202 (50%)	102 (100%)	0	100	100
3	C	93/192 (48%)	93 (100%)	0	100	100
3	F	93/192 (48%)	93 (100%)	0	100	100
3	I	93/192 (48%)	93 (100%)	0	100	100
All	All	3642/8070 (45%)	3635 (100%)	7 (0%)	93	96

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	185	CYS
1	J	600	LEU
1	K	477	THR
1	K	670	HIS
1	L	383	CYS
1	L	384	ASP
1	L	600	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	808	GLN
1	D	812	ASN

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 monosaccharides 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.

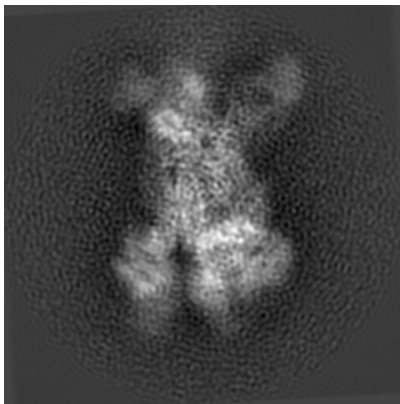
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8786. These allow visual inspection of the internal detail of the map and identification of artifacts.

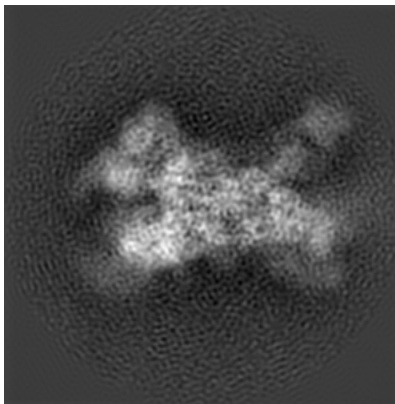
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

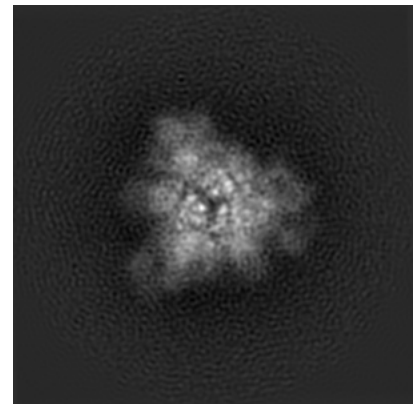
6.1.1 Primary map



X

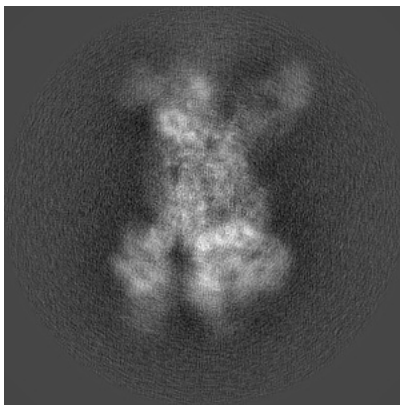


Y

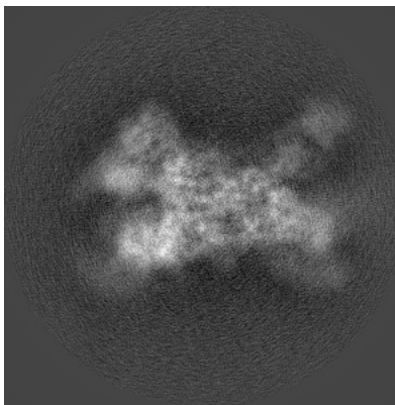


Z

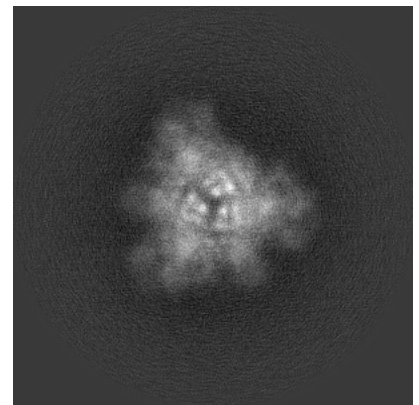
6.1.2 Raw map



X



Y

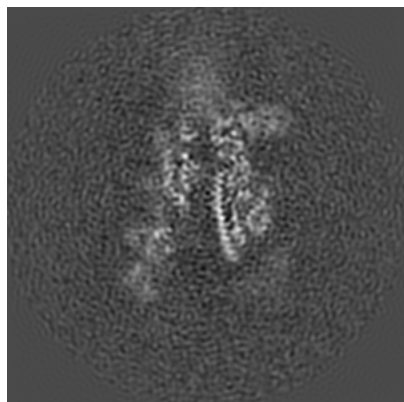


Z

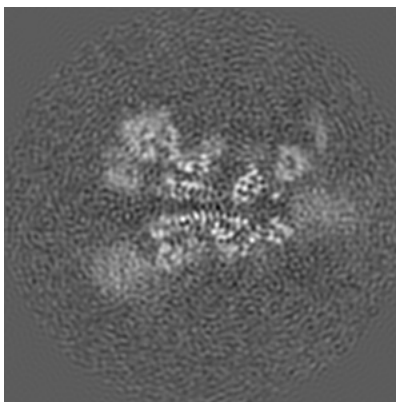
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

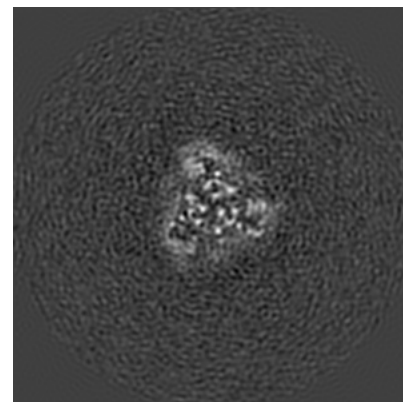
6.2.1 Primary map



X Index: 152

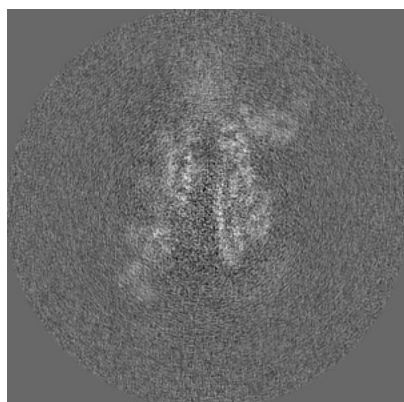


Y Index: 152

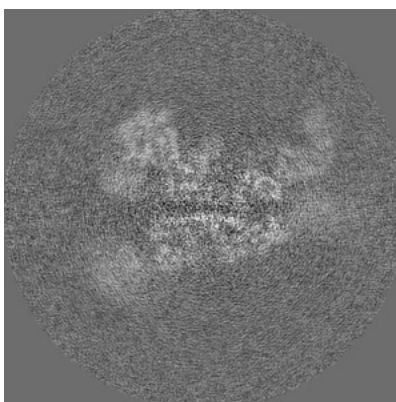


Z Index: 152

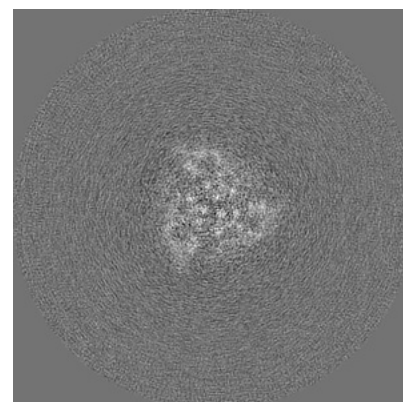
6.2.2 Raw map



X Index: 152



Y Index: 152

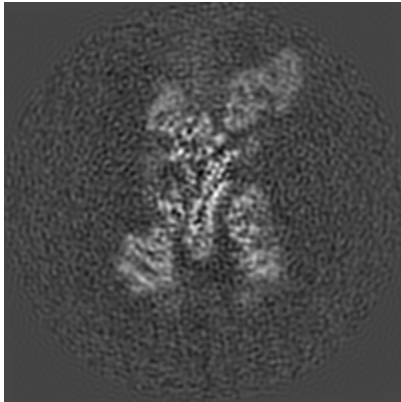


Z Index: 152

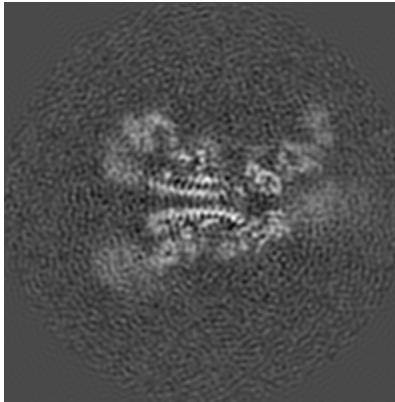
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

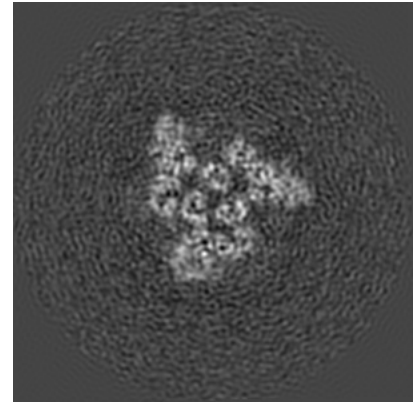
6.3.1 Primary map



X Index: 134

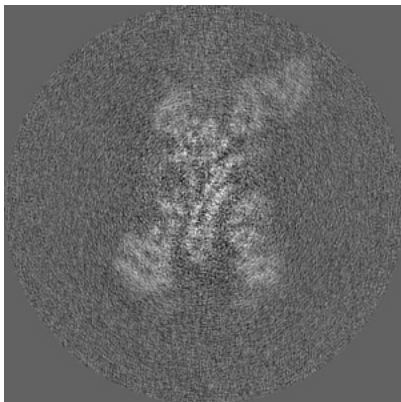


Y Index: 149

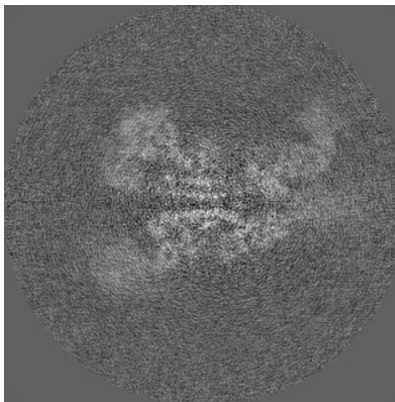


Z Index: 124

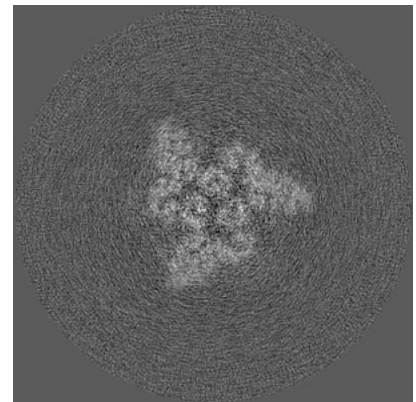
6.3.2 Raw map



X Index: 135



Y Index: 150

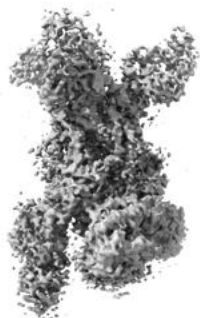


Z Index: 124

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

6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



X



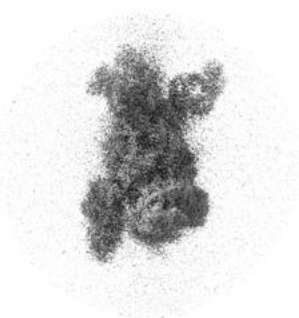
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.015. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

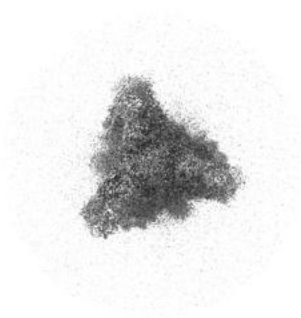
6.4.2 Raw map



X



Y



Z

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.

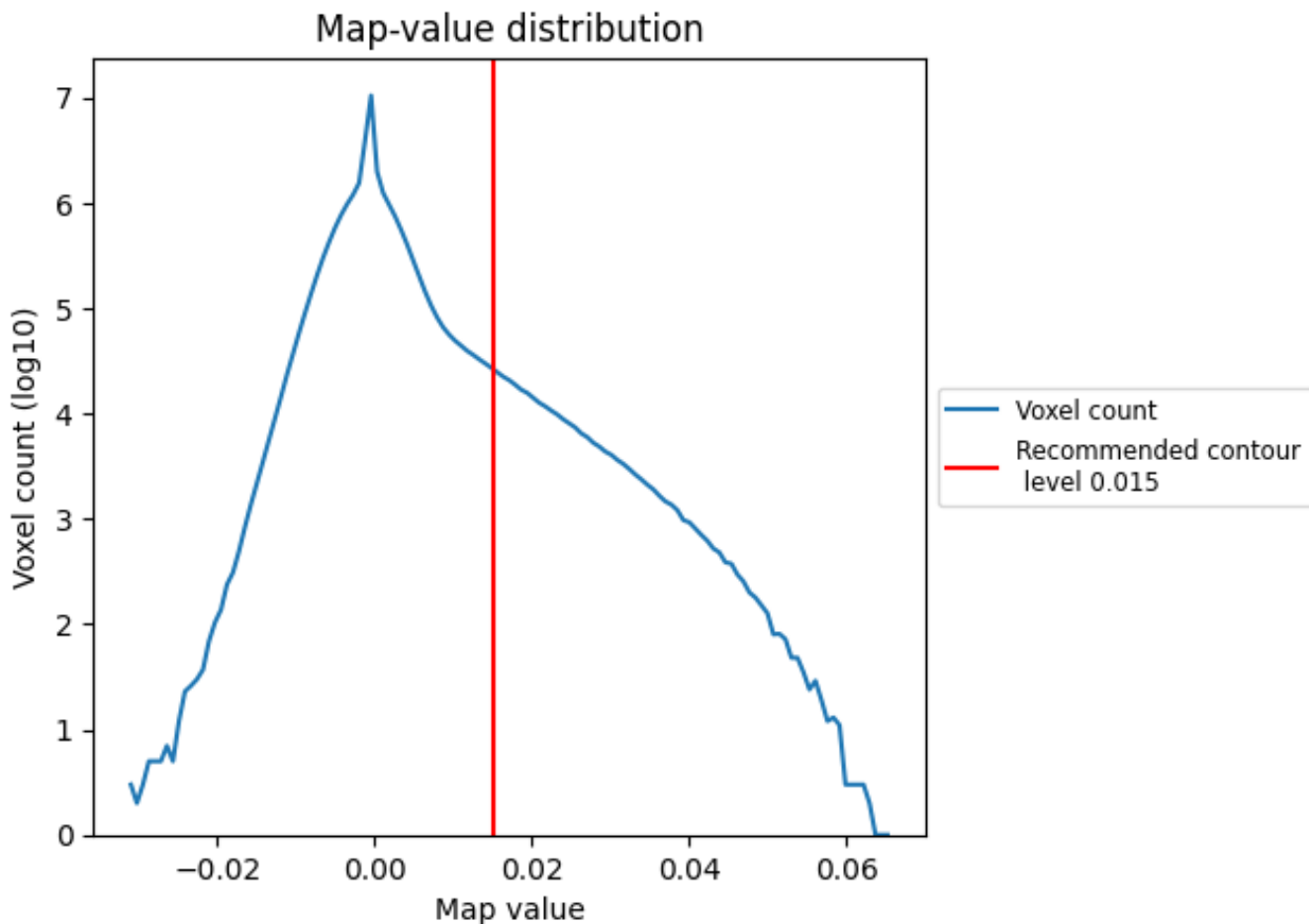
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

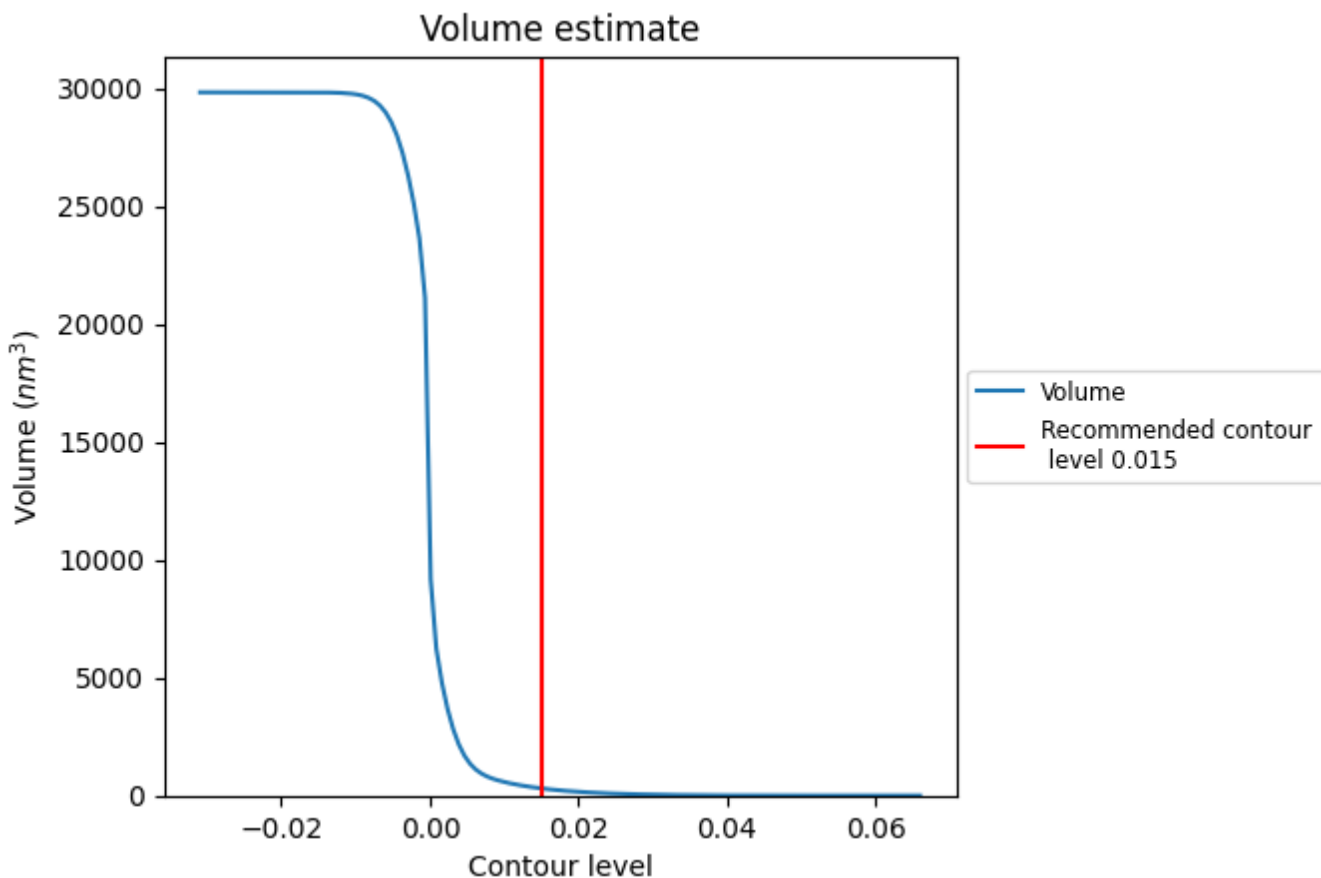
This section contains the results of statistical analysis of the map.

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.

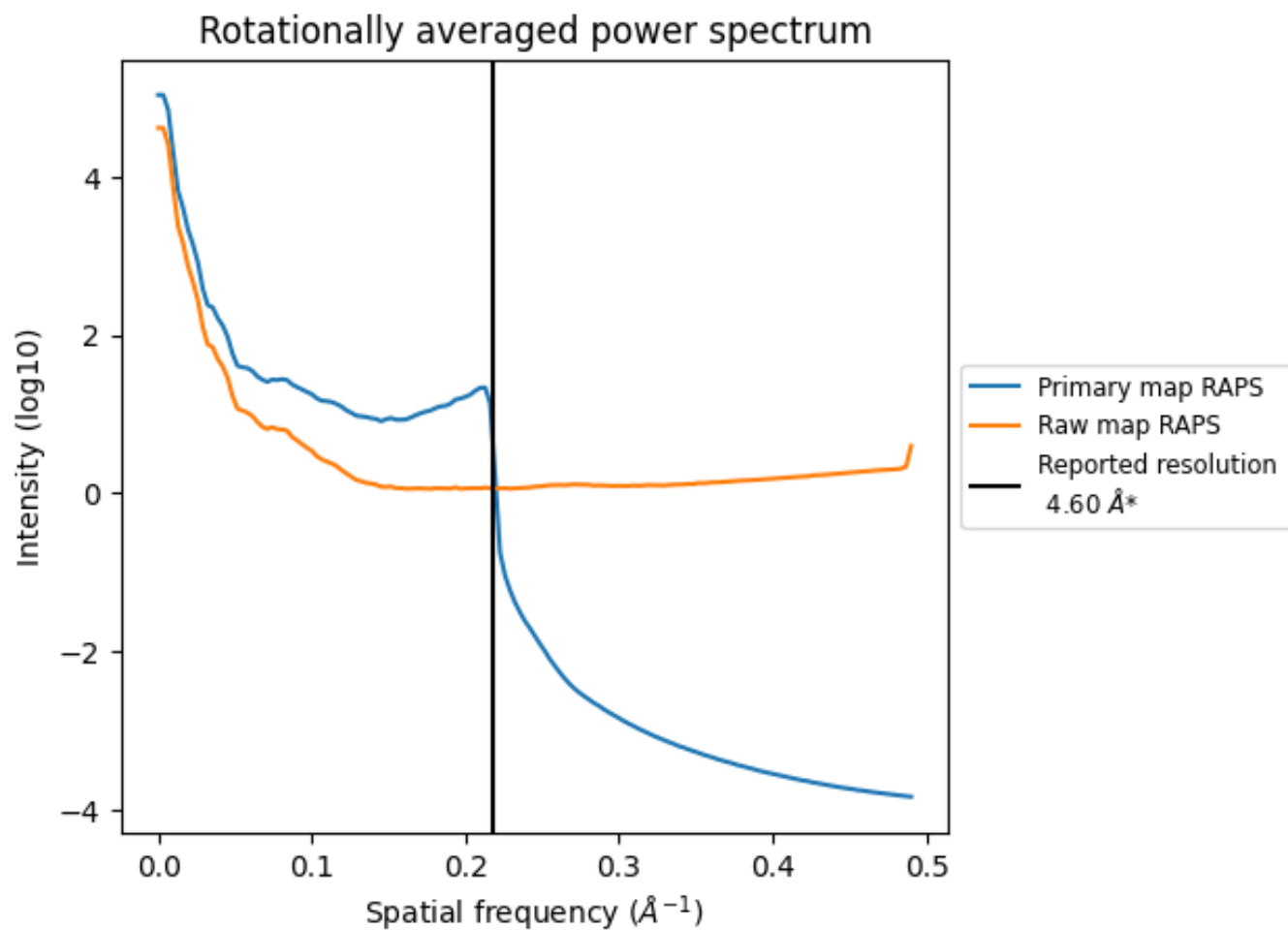
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 304 nm³; this corresponds to an approximate mass of 275 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 [i](#)

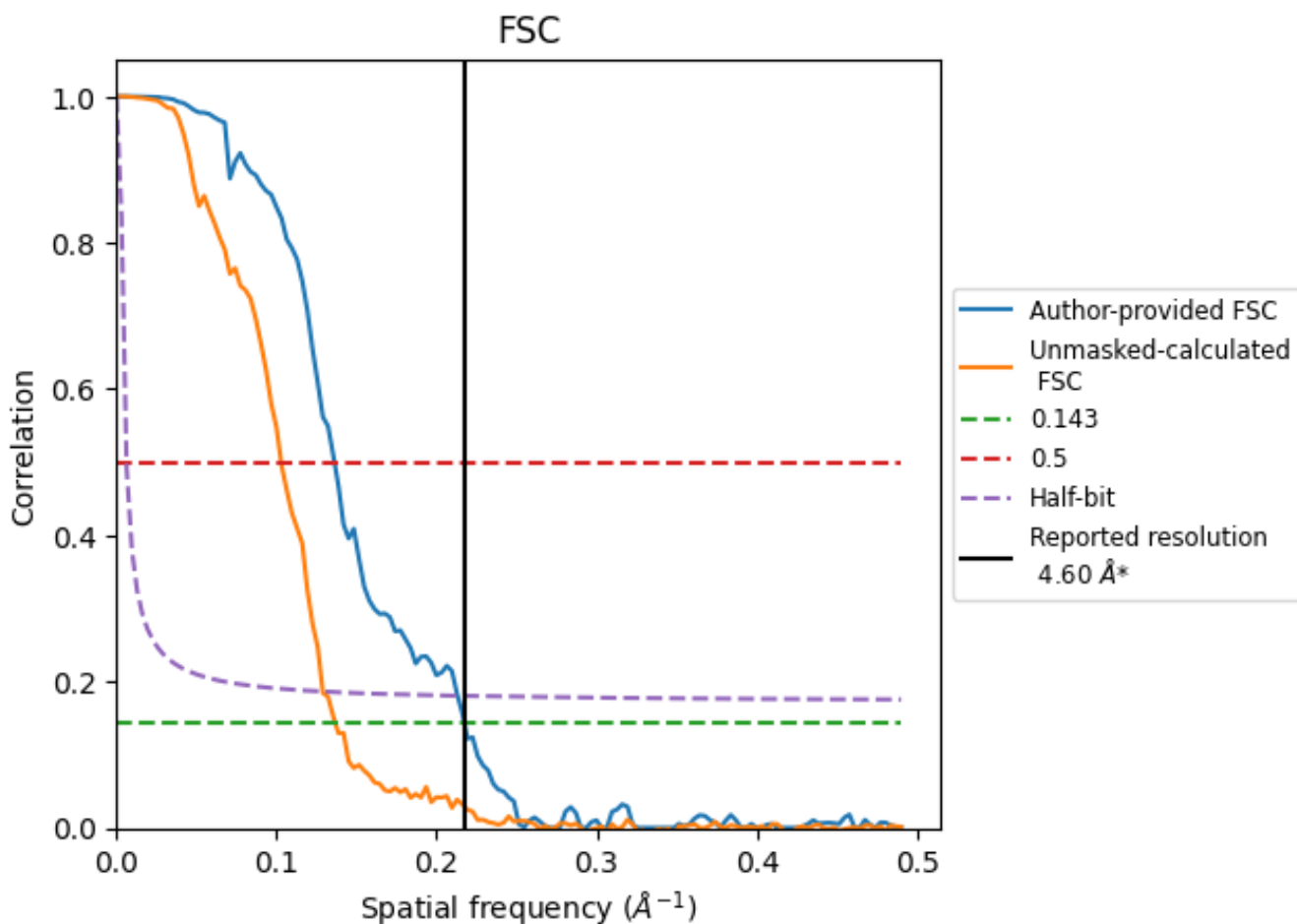


*Reported resolution corresponds to spatial frequency of 0.217 Å⁻¹

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 0.217 \AA^{-1}

8.2 Resolution estimates [i](#)

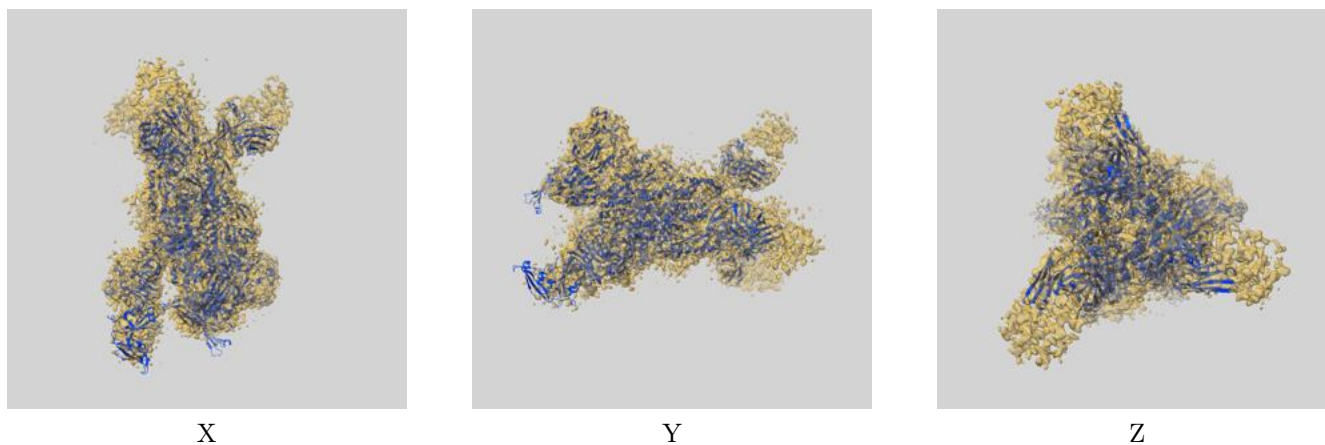
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.60	-	-
Author-provided FSC curve	4.60	7.34	4.69
Unmasked-calculated*	7.31	9.72	7.76

*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 7.31 differs from the reported value 4.6 by more than 10 %

9 Map-model fit [i](#)

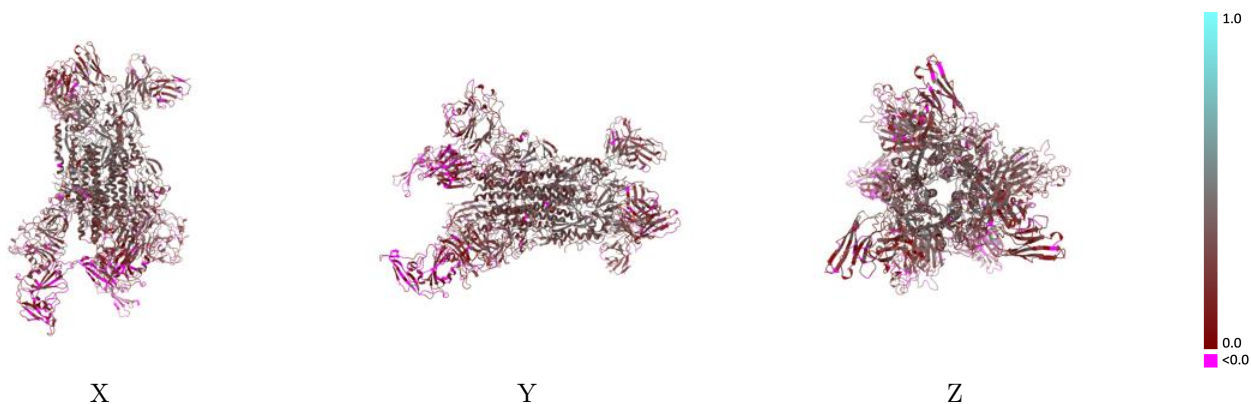
This section contains information regarding the fit between EMDB map EMD-8786 and PDB model 5W9K. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



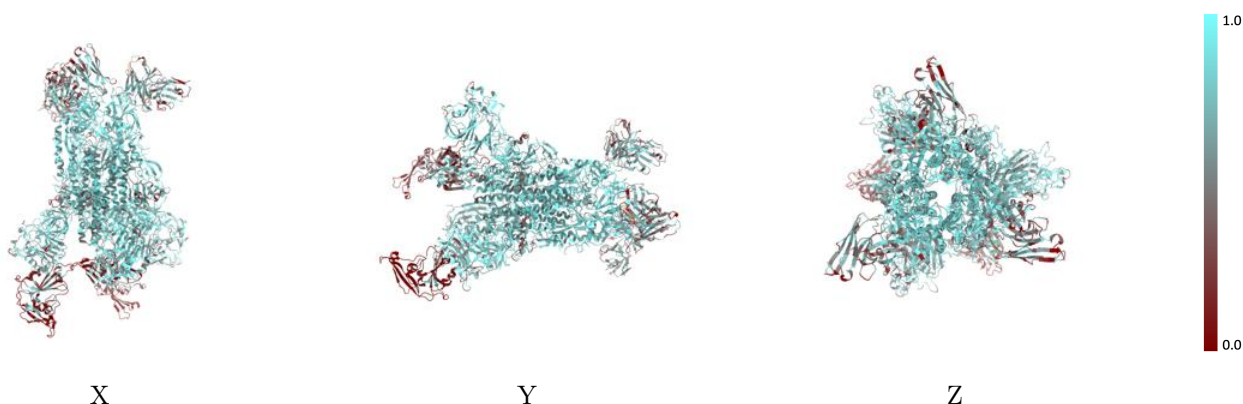
The images above show the 3D surface view of the map at the recommended contour level 0.015 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



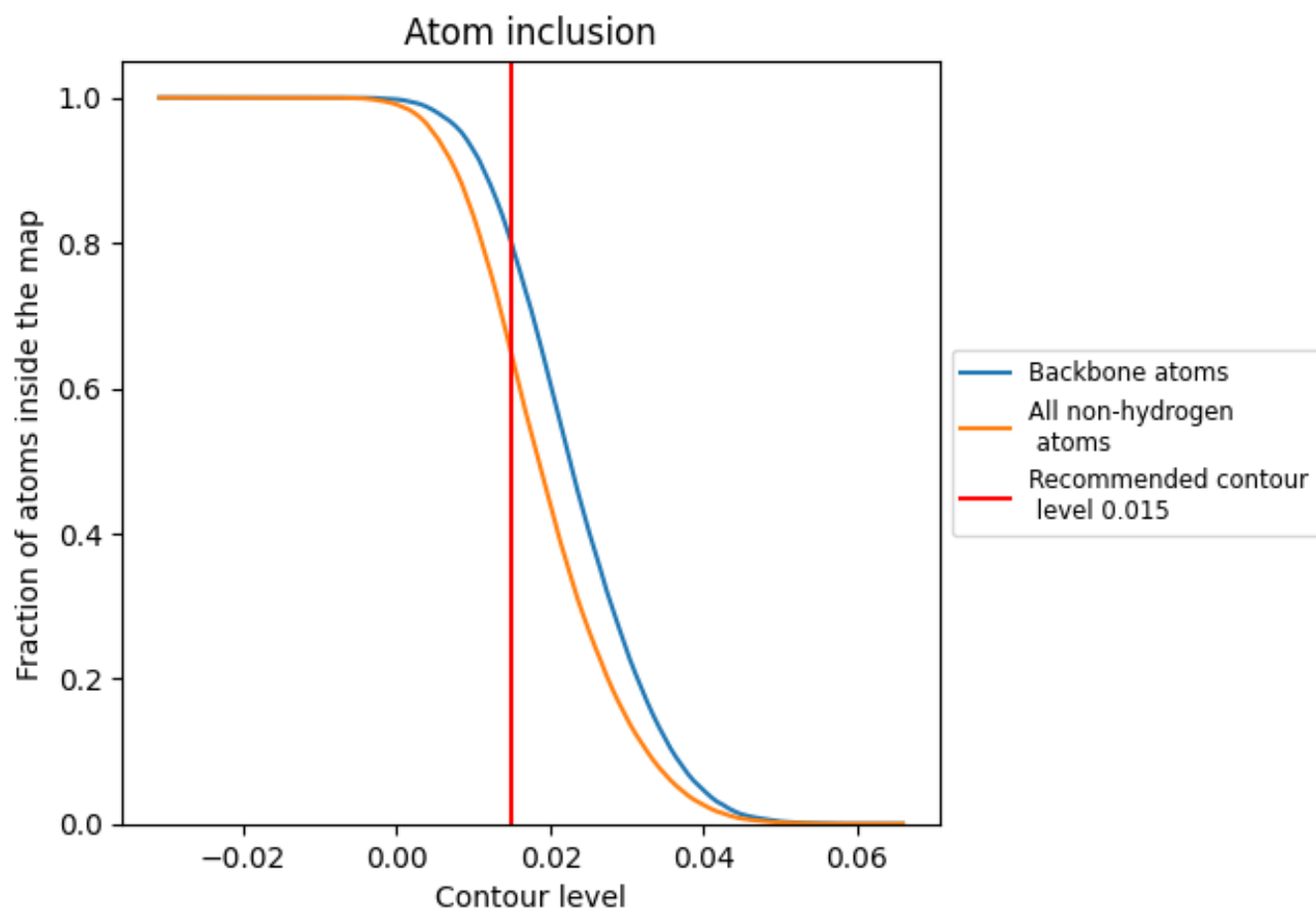
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.015).



























9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 64% 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 (0.015) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6444	 0.2420
A	 0.7549	 0.3120
B	 0.5535	 0.2290
C	 0.4660	 0.1920
D	 0.7657	 0.3240
E	 0.5686	 0.2430
F	 0.4697	 0.2050
G	 0.7626	 0.3280
H	 0.5665	 0.2480
I	 0.4442	 0.1870
J	 0.5622	 0.1840
K	 0.6602	 0.2120
L	 0.6137	 0.2020

