



wwPDB EM Validation Summary Report ⓘ

Jun 24, 2025 – 04:54 PM JST

PDB ID : 7X6A / pdb_00007x6a
EMDB ID : EMD-33019
Title : SARS-CoV-2 BA.2 variant spike protein in complex with Fab BD55-5840
Authors : Wang, X.; Wang, L.
Deposited on : 2022-03-07
Resolution : 3.50 Å(reported)

This is a wwPDB EM Validation Summary 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.dev118
MolProbity : 4-5-2 with Phenix2.0rc1
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

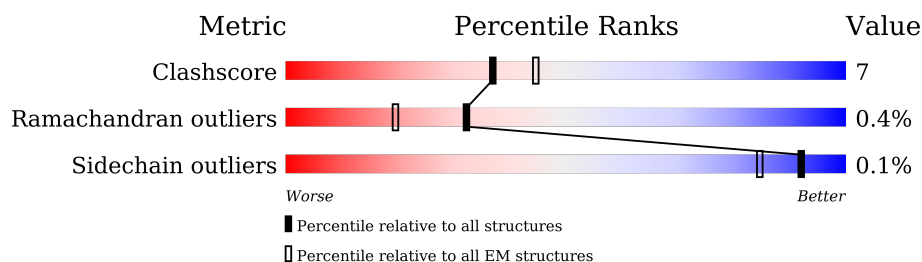
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1270	
1	B	1270	
1	D	1270	
2	G	119	
2	H	119	
2	I	119	
3	J	109	
3	K	109	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	L	109	<div><div></div><div>50%</div><div>76%</div><div>24%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 30879 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
1	A	1095	Total	C	N	O	S	0	0
			8548	5464	1424	1623	37		
1	B	1095	Total	C	N	O	S	0	0
			8548	5466	1423	1622	37		
1	D	1095	Total	C	N	O	S	0	0
			8551	5467	1424	1623	37		

There are 117 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ILE	THR	variant	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	27	SER	ALA	variant	UNP P0DTC2
A	142	ASP	GLY	variant	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	493	ARG	GLN	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	HIS	PRO	variant	UNP P0DTC2
A	683	ALA	ARG	variant	UNP P0DTC2
A	685	ALA	ARG	variant	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	variant	UNP P0DTC2
A	892	PRO	ALA	variant	UNP P0DTC2
A	899	PRO	ALA	variant	UNP P0DTC2
A	942	PRO	ALA	variant	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	22	ILE	THR	variant	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	27	SER	ALA	variant	UNP P0DTC2
B	142	ASP	GLY	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	493	ARG	GLN	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	683	ALA	ARG	variant	UNP P0DTC2
B	685	ALA	ARG	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	variant	UNP P0DTC2
B	892	PRO	ALA	variant	UNP P0DTC2
B	899	PRO	ALA	variant	UNP P0DTC2
B	942	PRO	ALA	variant	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
D	22	ILE	THR	variant	UNP P0DTC2
D	?	-	LEU	deletion	UNP P0DTC2
D	?	-	PRO	deletion	UNP P0DTC2
D	?	-	PRO	deletion	UNP P0DTC2
D	27	SER	ALA	variant	UNP P0DTC2
D	142	ASP	GLY	variant	UNP P0DTC2
D	213	GLY	VAL	variant	UNP P0DTC2
D	339	ASP	GLY	variant	UNP P0DTC2
D	371	PHE	SER	variant	UNP P0DTC2
D	373	PRO	SER	variant	UNP P0DTC2
D	375	PHE	SER	variant	UNP P0DTC2
D	376	ALA	THR	variant	UNP P0DTC2
D	405	ASN	ASP	variant	UNP P0DTC2
D	408	SER	ARG	variant	UNP P0DTC2
D	417	ASN	LYS	variant	UNP P0DTC2
D	440	LYS	ASN	variant	UNP P0DTC2
D	477	ASN	SER	variant	UNP P0DTC2
D	478	LYS	THR	variant	UNP P0DTC2
D	484	ALA	GLU	variant	UNP P0DTC2
D	493	ARG	GLN	variant	UNP P0DTC2
D	498	ARG	GLN	variant	UNP P0DTC2
D	501	TYR	ASN	variant	UNP P0DTC2
D	505	HIS	TYR	variant	UNP P0DTC2
D	614	GLY	ASP	variant	UNP P0DTC2
D	655	TYR	HIS	variant	UNP P0DTC2
D	679	LYS	ASN	variant	UNP P0DTC2
D	681	HIS	PRO	variant	UNP P0DTC2
D	683	ALA	ARG	variant	UNP P0DTC2
D	685	ALA	ARG	variant	UNP P0DTC2
D	764	LYS	ASN	variant	UNP P0DTC2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	796	TYR	ASP	variant	UNP P0DTC2
D	817	PRO	PHE	variant	UNP P0DTC2
D	892	PRO	ALA	variant	UNP P0DTC2
D	899	PRO	ALA	variant	UNP P0DTC2
D	942	PRO	ALA	variant	UNP P0DTC2
D	954	HIS	GLN	variant	UNP P0DTC2
D	969	LYS	ASN	variant	UNP P0DTC2
D	986	PRO	LYS	engineered mutation	UNP P0DTC2
D	987	PRO	VAL	engineered mutation	UNP P0DTC2

- Molecule 2 is a protein called Heavy chain of Fab BD55-5840.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	119	Total	C	N	O	S	0	0
			927	585	157	182	3		
2	G	119	Total	C	N	O	S	0	0
			927	585	157	182	3		
2	I	119	Total	C	N	O	S	0	0
			927	585	157	182	3		

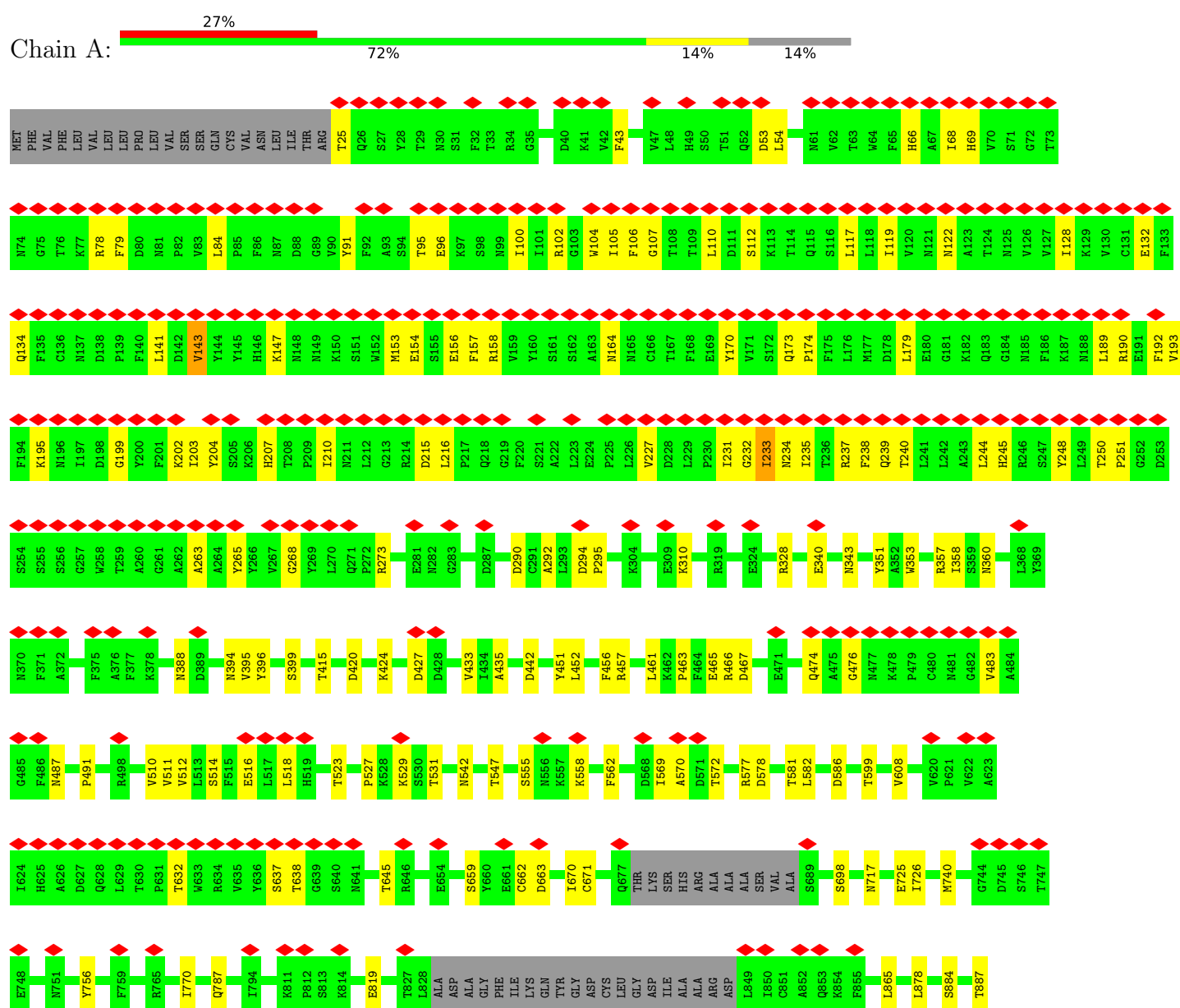
- Molecule 3 is a protein called Light chain of Fab BD55-5840.

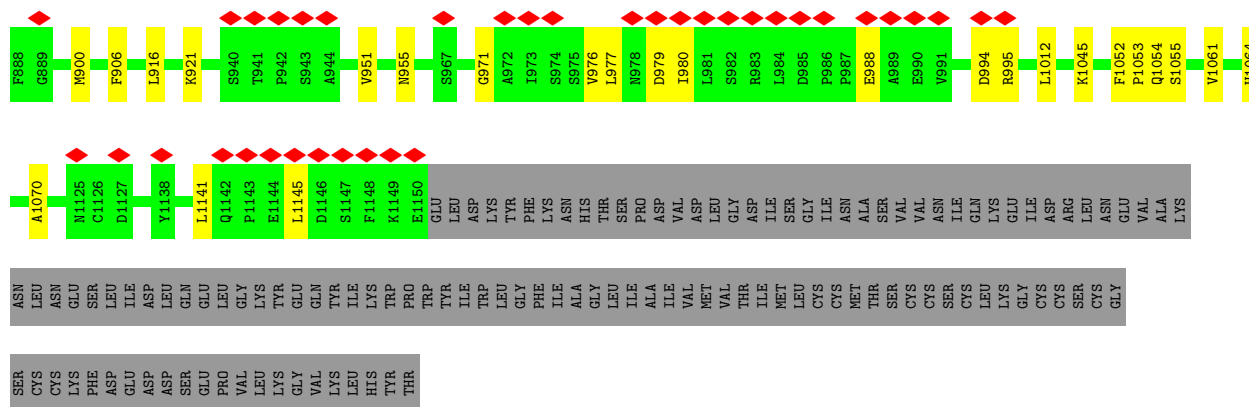
Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	109	Total	C	N	O	S	0	0
			817	510	142	162	3		
3	J	109	Total	C	N	O	S	0	0
			817	510	142	162	3		
3	K	109	Total	C	N	O	S	0	0
			817	510	142	162	3		

3 Residue-property plots

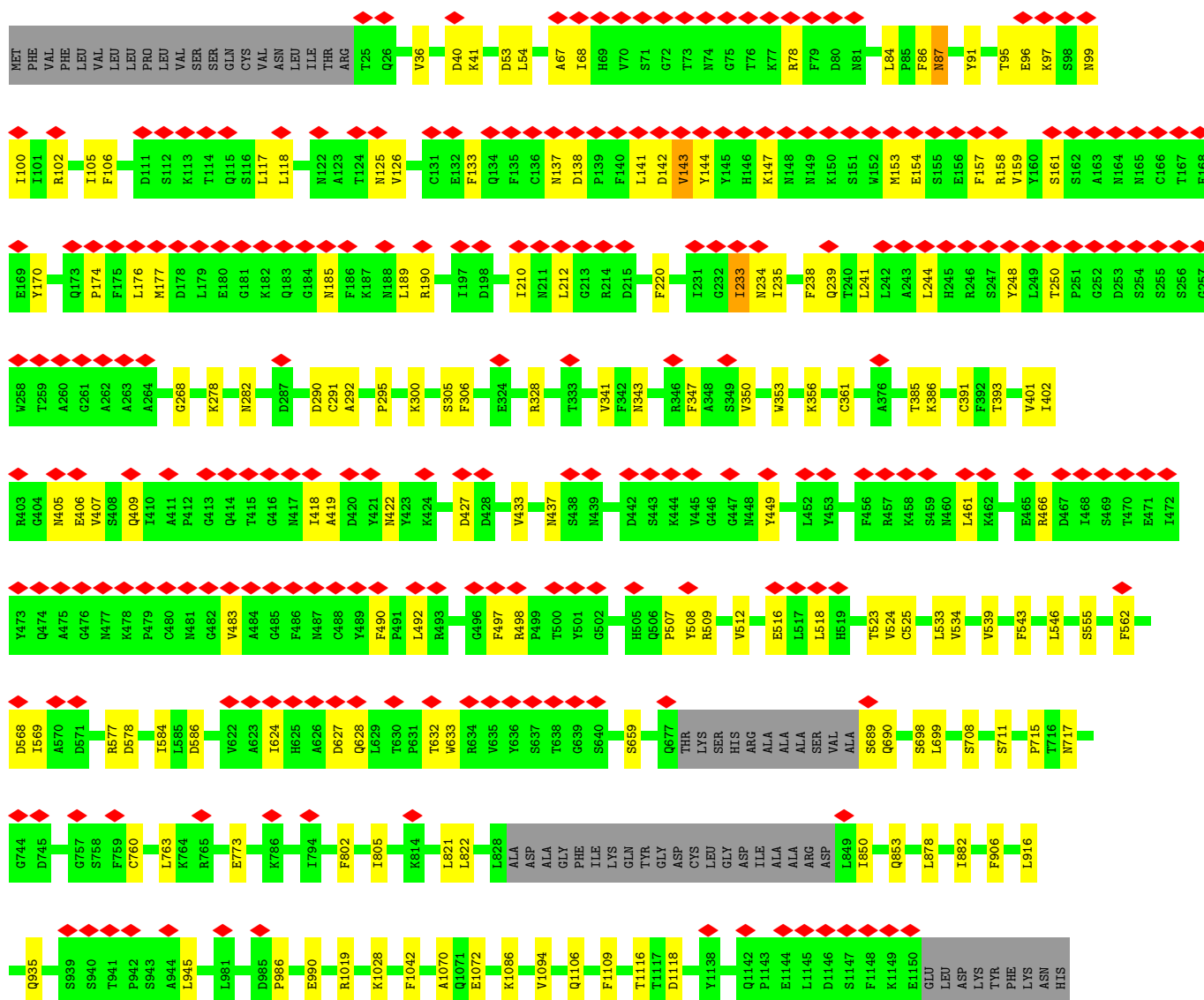
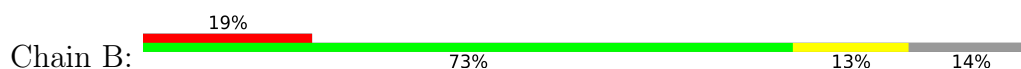
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Spike glycoprotein



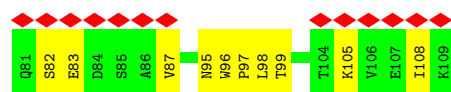


• Molecule 1: Spike glycoprotein

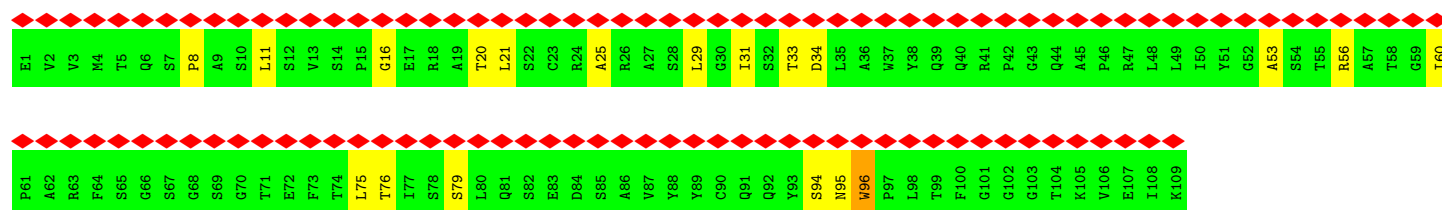
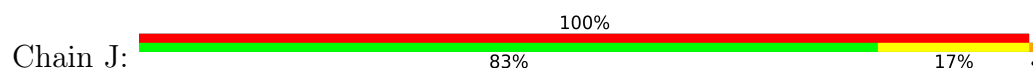


- Molecule 1: Spike glycoprotein

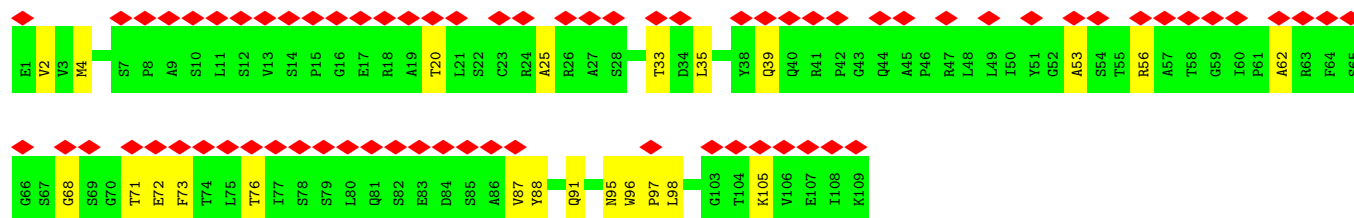
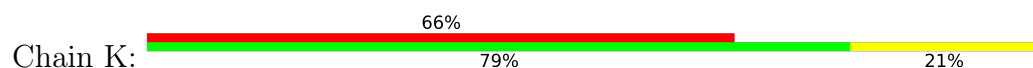
P812	P631	T553	L492	C432	A372	K304	F220	S155	Y91	MET
S813	T632	E554	R493	V433	P373	S305	L223	E156	S94	PHE
K814	W633	S494	S494	I434	F374	F306		E157	T95	PHE
E819	R634	N556	Y495	A435	F375			R158	E96	LEU
L828	W636	K558	Q496	W436	F377	N317	L226	V159	K97	VAL
ALA	S637	F559	R498	R438	K378	F318	V227	Y160	S98	LEU
ASP	T638	L560	P499	W439	C379	V320	D228	Y160	S97	LEU
GLY	ALA	P561	T500	K440	Y380	R319	L229	S161	N99	PRO
PHE	S639	F562	Y501	L441	G381	Q321	P230	S162	I100	LEU
S640	ILE	Q563	Q563	D442	W382	T323	T231	A163	I101	VAL
ILE	W641	Q564	Q564	S443	S383	E324	G232	N164	I102	SER
LYS	W642	Q564	W503	S443	S383	S325	G233	M165	R102	SER
GLN		F565	Q504	K444	P384	S326	W234	C166	I105	GLN
TYR	L650	F566	H505	V445	T385	V327	T235	C167	F106	CYS
GLY	L651	R567	H505	V445	T385	V327	T236	F168	G107	ASN
ASP		D568	Q506	D446	K386	R328	R237	E169	F106	VAL
CYS	S659	L569	P507	Q447	K386	F329	F238		G107	LEU
LEU		A570	Y508	W448	L387	P330	Q239		T108	THR
GLY	T676	A570	R509	Y449	L387	N331	T240		T109	
ASP	Q677	W572	W510	W450	L390	S332	L241	Q173	L110	ARG
ILE	THR	L572	W511	Y451	C391	T333	L242	P174	D111	T25
ALA	LYS	D574	W512	L452	F392	N334	A243	F175	S112	Q26
ALA	SER	L513	L513	Y453	T393	L335	A243	K113	K113	T33
ARG	ARG	A575	S514	R454	N394	C336	H245	T114	Q115	
ALA	ALA	W576	S515	L455	N394	F338	R246	D178	S116	V36
ALA	ALA	S577	F515	L455	Y395	P337	S247	E180	L117	Y37
ALA	SER	D578	E516	F456	Y396	F338	Y248	G181	L118	
VAL	P579	P579	L517	R457	A397	D339	L249	K182	I119	
ALA	Q580	Q580	L517	R457	A397	E340	T250	Q183	W120	
S689	T581	H519	H519	K458	D398	S341	P251	G184	N121	
S698	L582	A520	A520	S459	S399	F342	D253	N185	A123	D53
S698	E583	P521	T523	L461	V401	N343	S254	F186	T124	L54
E702	L584	A522	T523	K462	I402	A344	S254	K187		F65
N703	L585	L585	W524	F464	G404	R346	S256	N188	T128	H66
S708	D586	S587	C525	E465	N405	F347	S256	L189	K129	A67
S711	L588	T588	G526	R466	E406	A348	Q257	R190	W130	I68
Q935	P589	C590	P527	D467	W407	S349	W258	E191	G131	H69
T716	T716	S591	K528	E466	N405	A348	Q258	E191	E132	V70
W717	F592	K529	S529	S469	Q409	S351	T259	F192	F133	S71
T599	T599	S530	S530	T470	I410	A352	G261	K195	Q134	G72
W608	W608	T531	T531	E471	A411	K353	A262	N196	F135	T73
N616	N616	N532	N532	F471	A411	N354	A263	I197	C136	W74
C617	C617	L533	L533	Y473	G413	R355	A264	D198	D138	T76
T618	T618	W534	W534	Q474	Q414	K356	Y265	G199	W139	K77
E619	E619	W535	W535	Q474	Q414	R357	Y266	Y200	F140	R78
W620	W620	N536	N536	A475	T415	S357				



• Molecule 3: Light chain of Fab BD55-5840



• Molecule 3: Light chain of Fab BD55-5840



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	355445	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.043	Depositor
Minimum map value	-1.849	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.061	Depositor
Recommended contour level	0.558	Depositor
Map size (\AA)	385.2, 385.2, 385.2	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.07, 1.07, 1.07	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	0.11	0/8758	0.30	0/11930
1	B	0.11	0/8758	0.28	0/11930
1	D	0.12	0/8761	0.30	0/11934
2	G	0.08	0/952	0.23	0/1295
2	H	0.11	0/952	0.31	0/1295
2	I	0.10	0/952	0.27	0/1295
3	J	0.10	0/834	0.28	0/1133
3	K	0.10	0/834	0.28	0/1133
3	L	0.11	0/834	0.33	0/1133
All	All	0.11	0/31635	0.29	0/43078

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	J	0	1
3	K	0	1
3	L	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	J	96	TRP	Peptide
3	K	96	TRP	Peptide
3	L	96	TRP	Peptide

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	8548	0	8314	123	0
1	B	8548	0	8319	107	0
1	D	8551	0	8323	157	0
2	G	927	0	862	11	0
2	H	927	0	862	15	0
2	I	927	0	862	15	0
3	J	817	0	802	13	0
3	K	817	0	802	14	0
3	L	817	0	802	15	0
All	All	30879	0	29948	447	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 7.

The worst 5 of 447 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:D:185:ASN:HB2	1:D:212:LEU:CD2	1.38	1.50
1:B:185:ASN:O	1:B:212:LEU:HD23	1.36	1.24
1:D:185:ASN:CB	1:D:212:LEU:CD2	2.25	1.15
1:B:185:ASN:O	1:B:212:LEU:CD2	1.98	1.12
1:D:185:ASN:HB2	1:D:212:LEU:HD21	1.27	1.11

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	1089/1270 (86%)	1023 (94%)	62 (6%)	4 (0%)	30	64
1	B	1089/1270 (86%)	1024 (94%)	60 (6%)	5 (0%)	25	59
1	D	1089/1270 (86%)	1007 (92%)	77 (7%)	5 (0%)	25	59
2	G	117/119 (98%)	116 (99%)	1 (1%)	0	100	100
2	H	117/119 (98%)	111 (95%)	6 (5%)	0	100	100
2	I	117/119 (98%)	113 (97%)	4 (3%)	0	100	100
3	J	107/109 (98%)	96 (90%)	11 (10%)	0	100	100
3	K	107/109 (98%)	97 (91%)	10 (9%)	0	100	100
3	L	107/109 (98%)	95 (89%)	12 (11%)	0	100	100
All	All	3939/4494 (88%)	3682 (94%)	243 (6%)	14 (0%)	32	64

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	ILE
1	B	143	VAL
1	B	233	ILE
1	D	143	VAL
1	D	233	ILE

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	950/1109 (86%)	950 (100%)	0	100	100
1	B	950/1109 (86%)	949 (100%)	1 (0%)	92	97
1	D	951/1109 (86%)	950 (100%)	1 (0%)	92	97
2	G	98/98 (100%)	98 (100%)	0	100	100
2	H	98/98 (100%)	98 (100%)	0	100	100
2	I	98/98 (100%)	98 (100%)	0	100	100
3	J	88/88 (100%)	88 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	K	88/88 (100%)	88 (100%)	0	100	100
3	L	88/88 (100%)	88 (100%)	0	100	100
All	All	3409/3885 (88%)	3407 (100%)	2 (0%)	92	97

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	87	ASN
1	D	87	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	82	GLN
2	G	3	GLN
2	I	43	GLN
1	B	1106	GLN
1	B	965	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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.

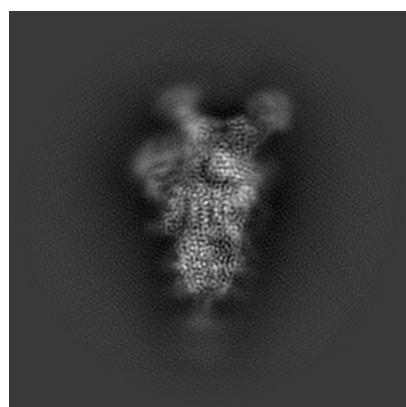
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33019. These allow visual inspection of the internal detail of the map and identification of artifacts.

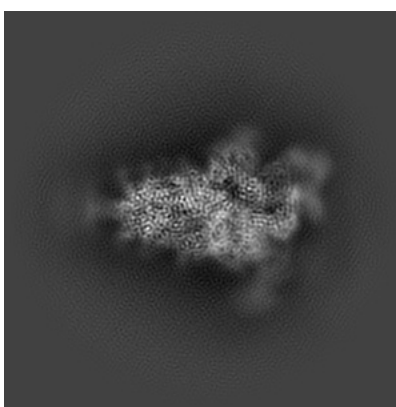
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

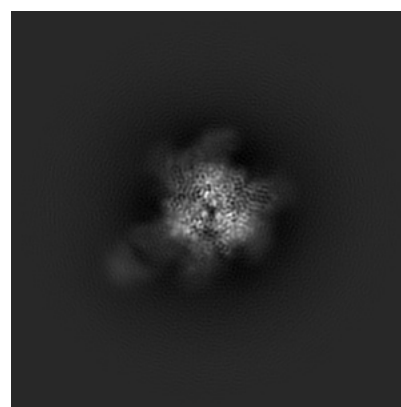
6.1.1 Primary map



X



Y

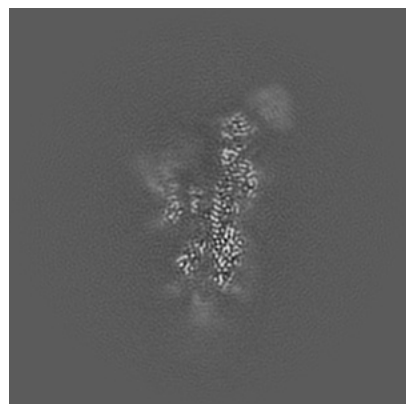


Z

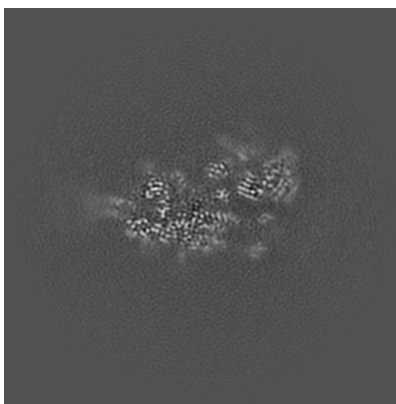
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

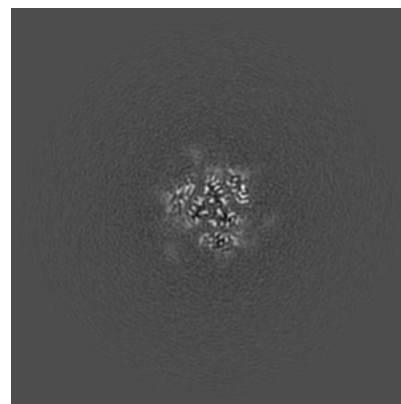
6.2.1 Primary map



X Index: 180



Y Index: 180

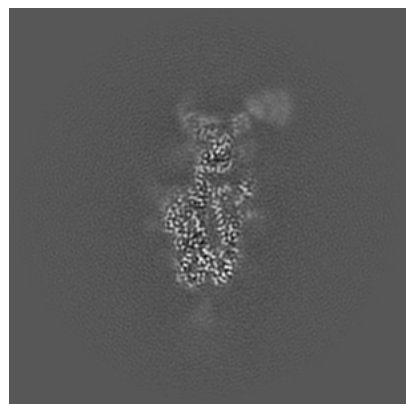


Z Index: 180

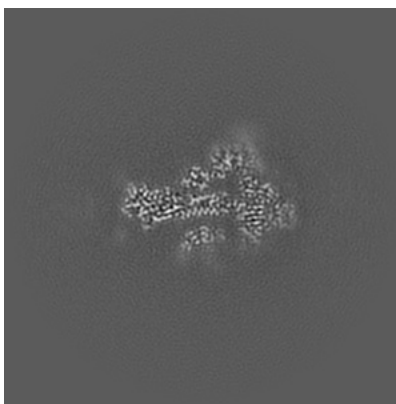
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

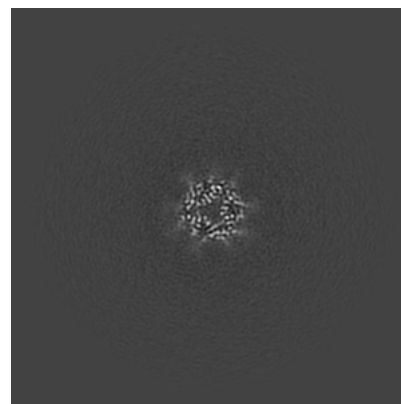
6.3.1 Primary map



X Index: 192



Y Index: 195



Z Index: 129

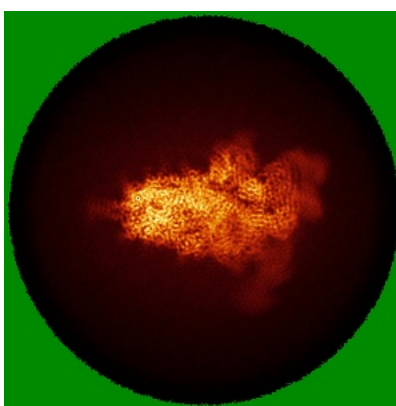
The images above show the largest variance slices of the map in three orthogonal directions.

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

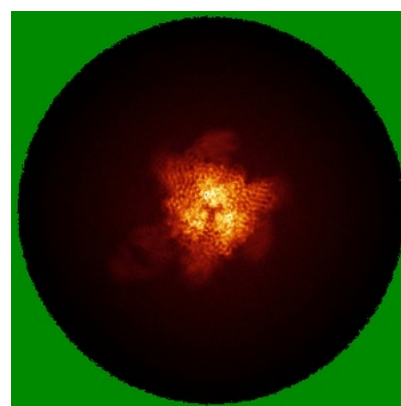
6.4.1 Primary map



X



Y



Z

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 0.558. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

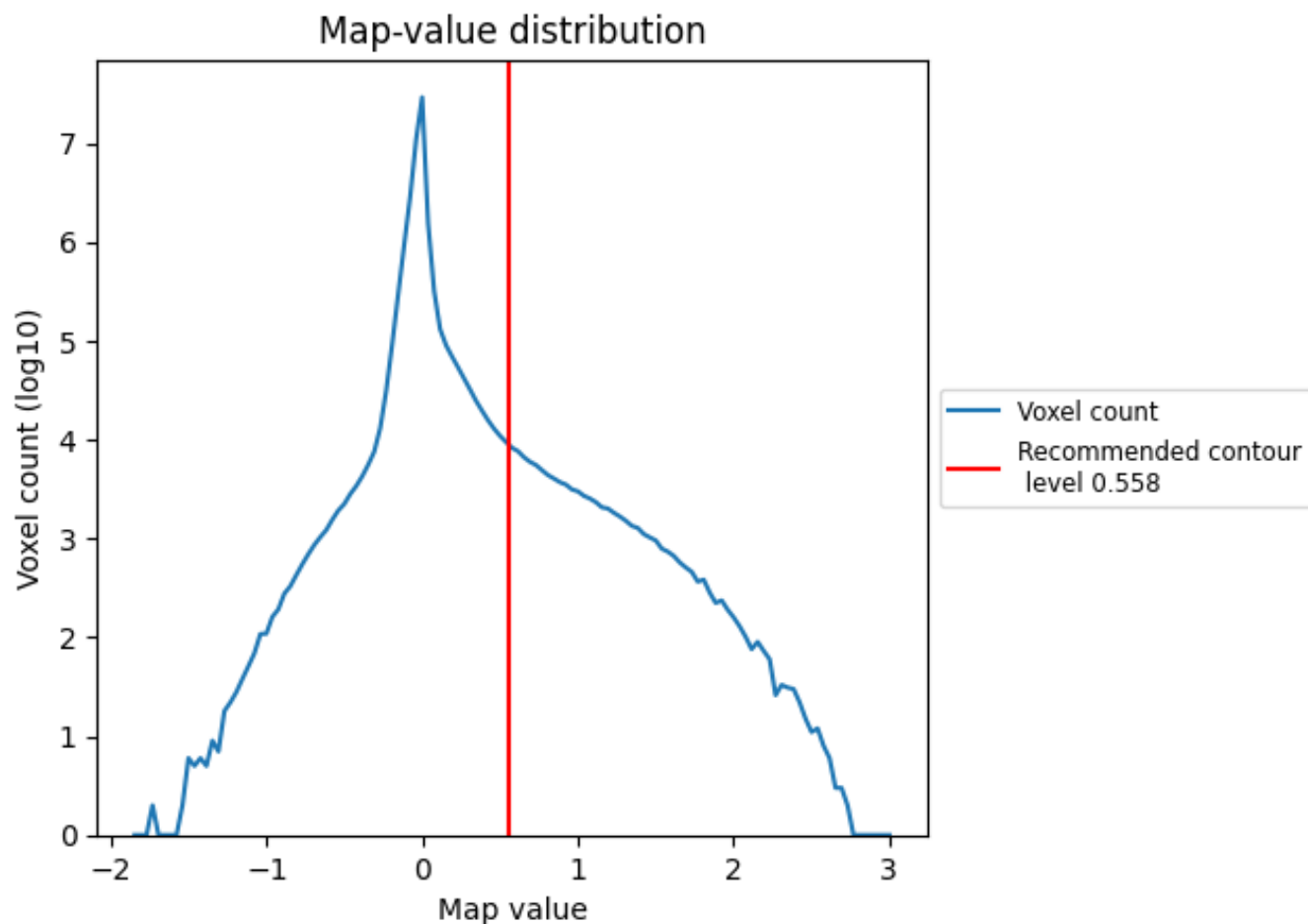
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

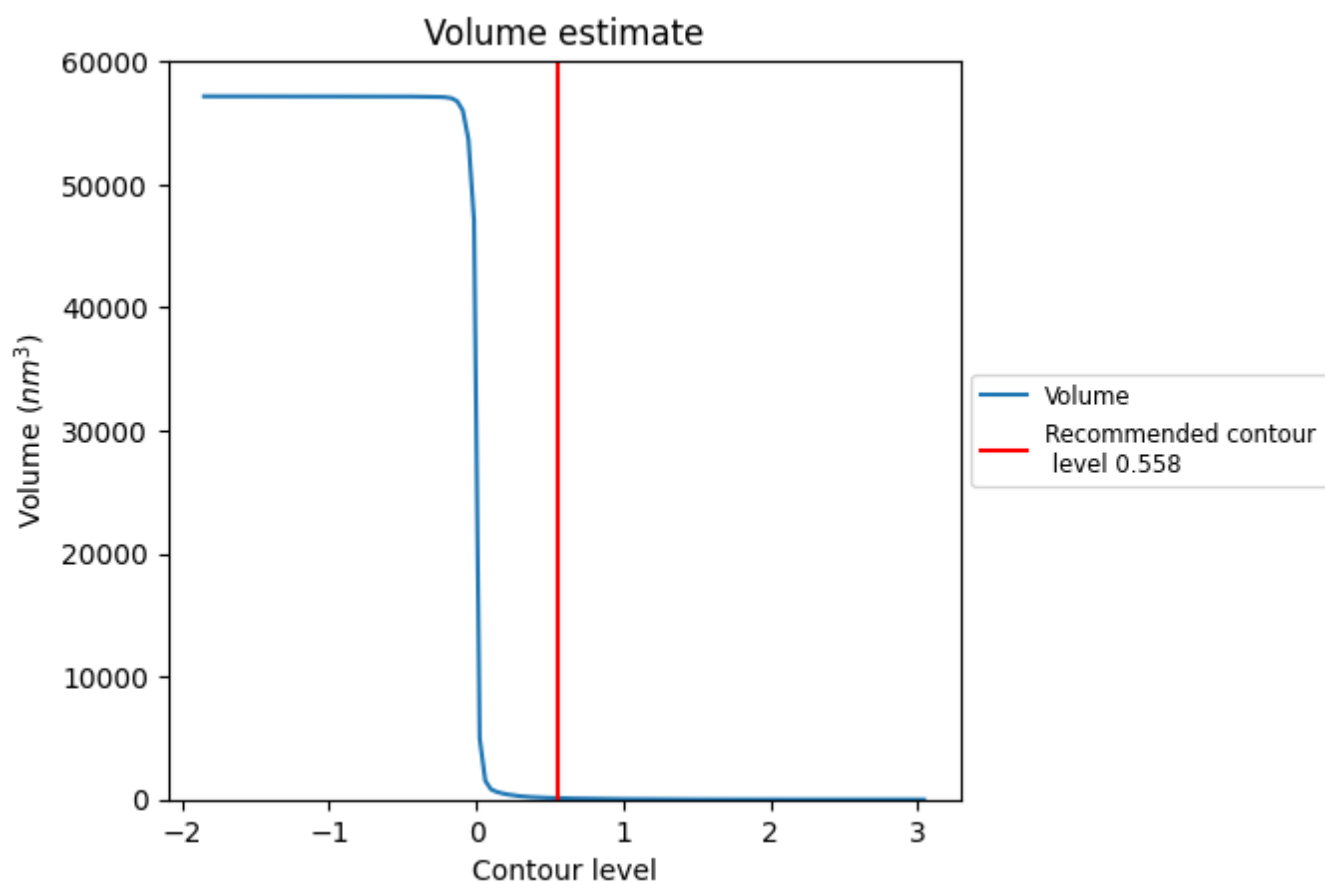
This section contains the results of statistical analysis of the map.

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.

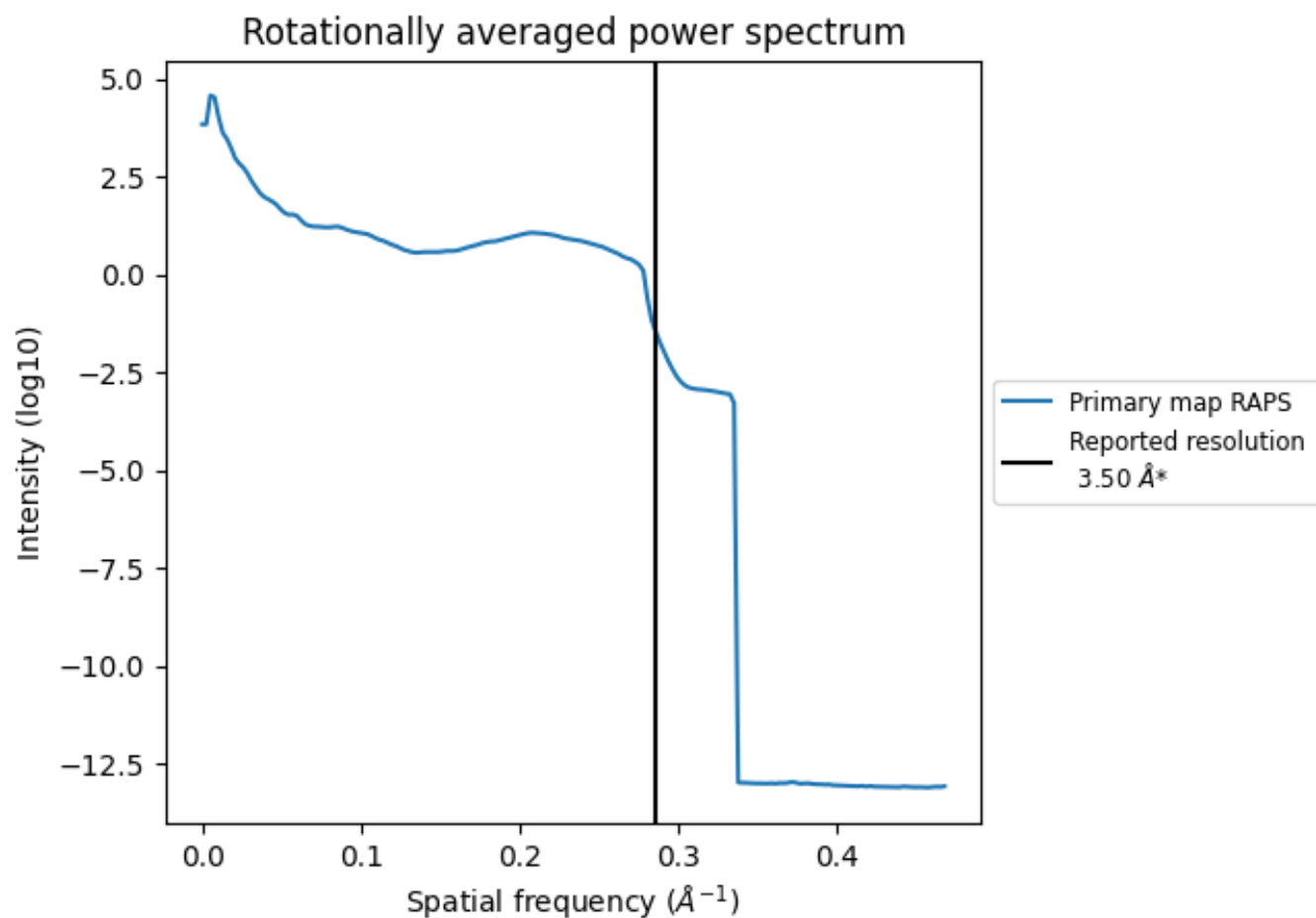
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 115 nm³; this corresponds to an approximate mass of 104 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 0.286 \AA^{-1}

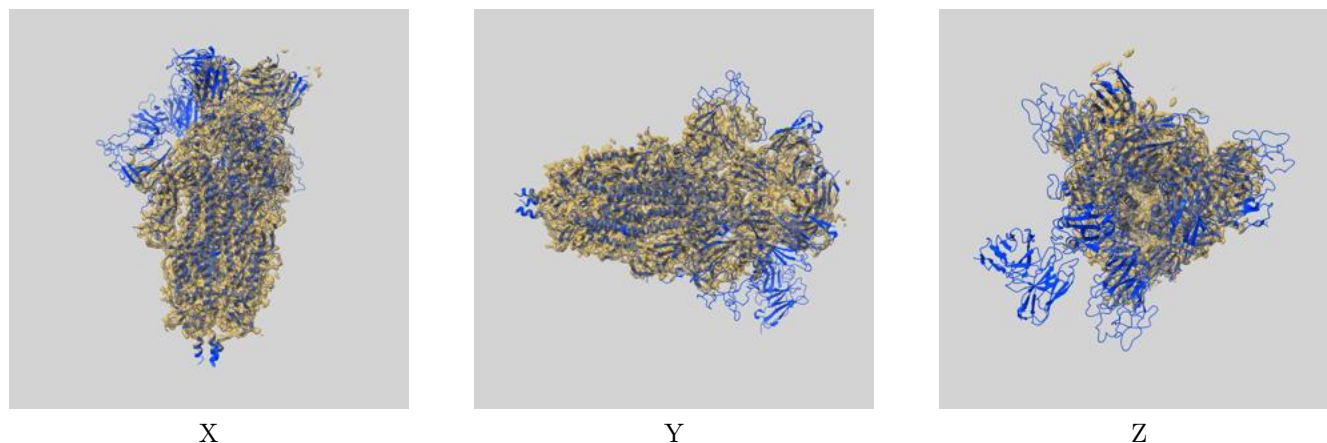
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

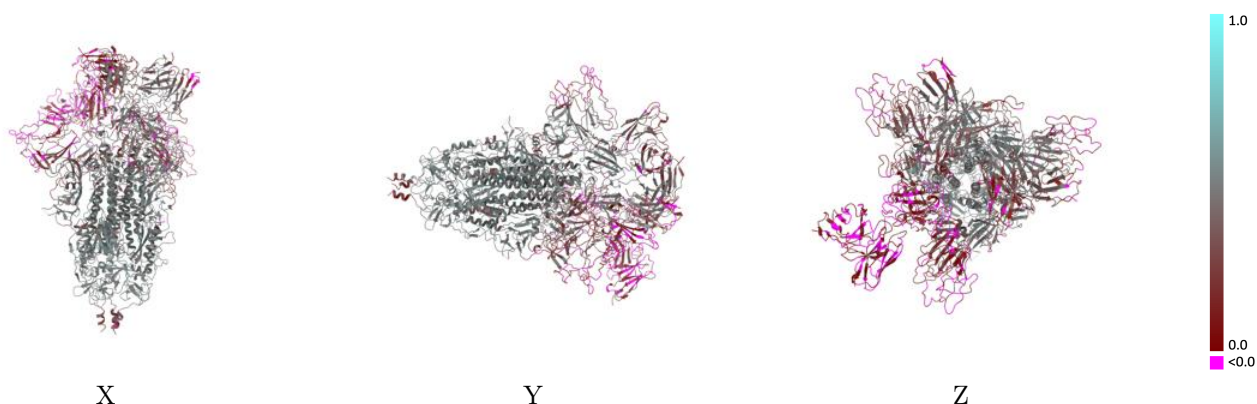
This section contains information regarding the fit between EMDB map EMD-33019 and PDB model 7X6A. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



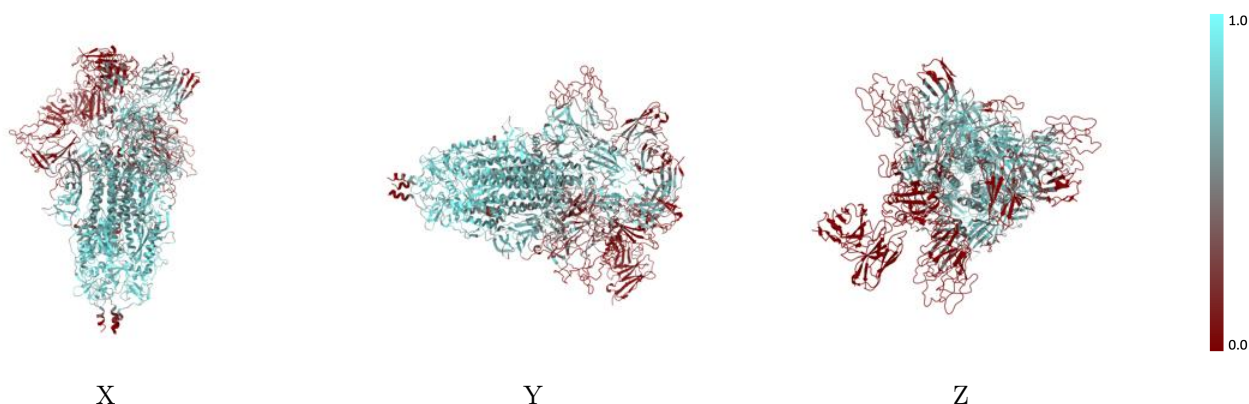
The images above show the 3D surface view of the map at the recommended contour level 0.558 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



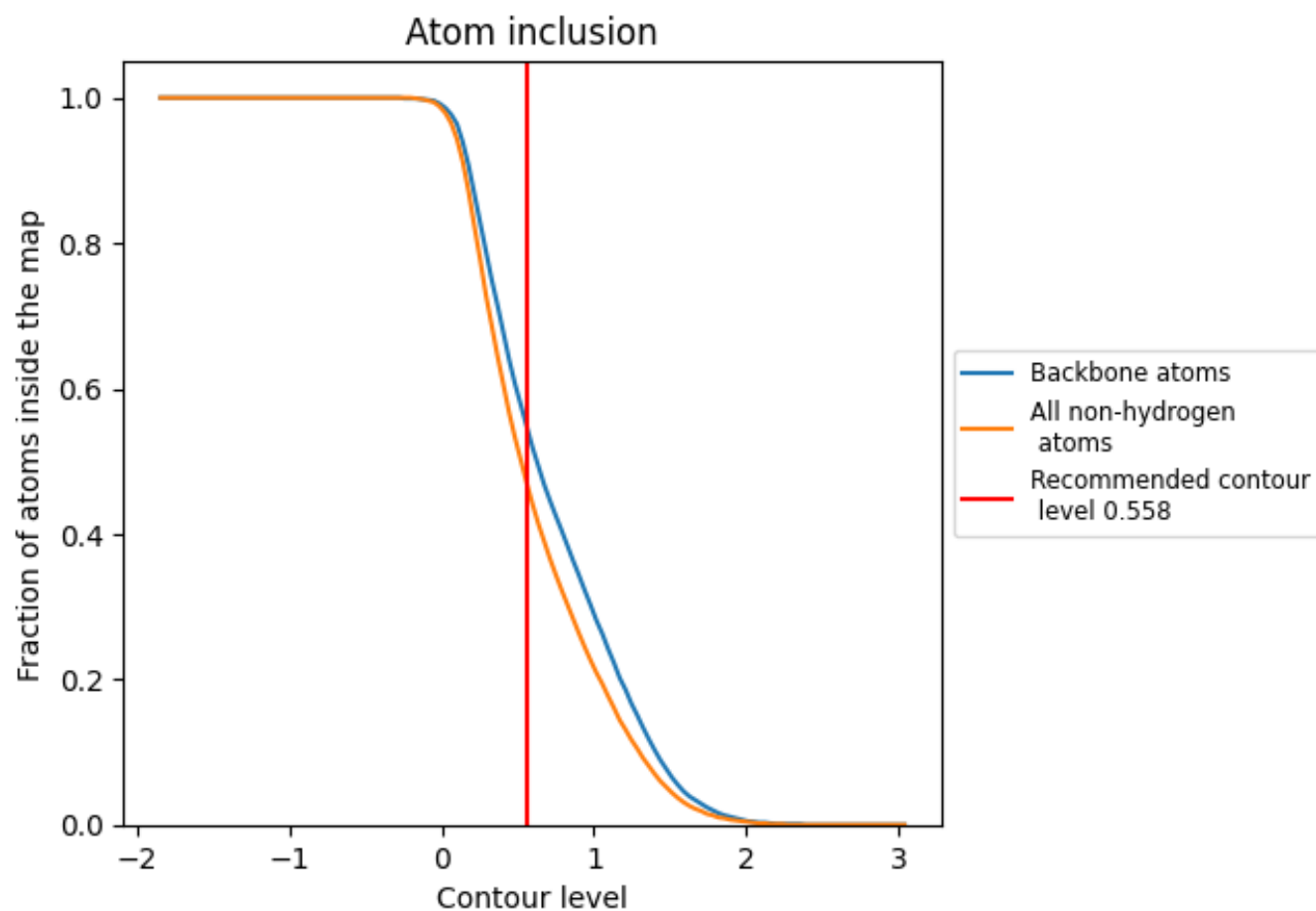
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.558).

9.4 Atom inclusion [i](#)



At the recommended contour level, 55% of all backbone atoms, 47% 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 (0.558) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.4710	<div></div> 0.3560
A	<div></div> 0.5280	<div></div> 0.3920
B	<div></div> 0.5960	<div></div> 0.4180
D	<div></div> 0.4380	<div></div> 0.3340
G	<div></div> 0.0000	<div></div> 0.0540
H	<div></div> 0.4690	<div></div> 0.3850
I	<div></div> 0.1870	<div></div> 0.2420
J	<div></div> 0.0000	<div></div> 0.0480
K	<div></div> 0.2860	<div></div> 0.3250
L	<div></div> 0.4060	<div></div> 0.3150

1.0

0.0

<0.0