



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 07:18 PM UTC

PDB ID : 8VGW / pdb_00008vgw
EMDB ID : EMD-43228
Title : VRC01 Fab bound to the HIV-1 CH848 DE3 SOSIP
Authors : Henderson, R.; Acharya, P.
Deposited on : 2023-12-29
Resolution : 3.90 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:

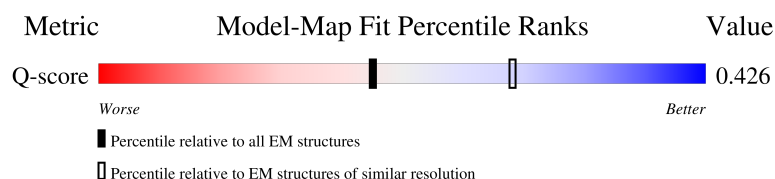
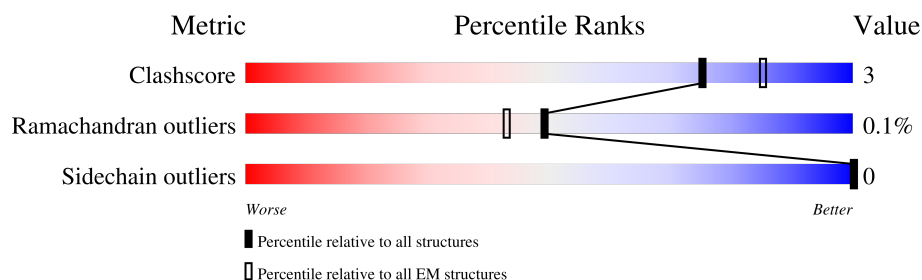
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.90 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.



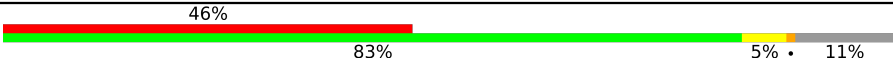
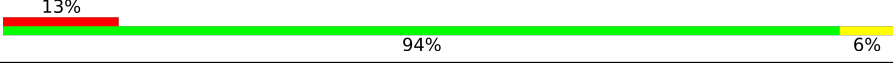
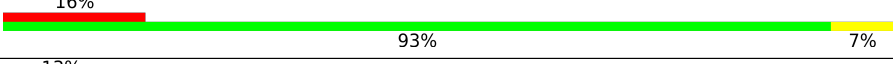
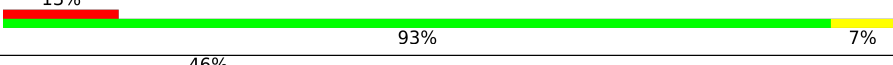
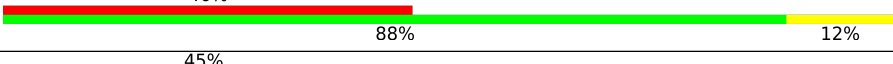
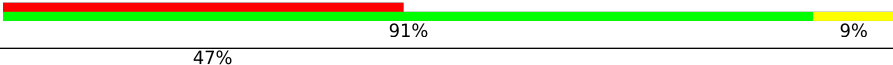
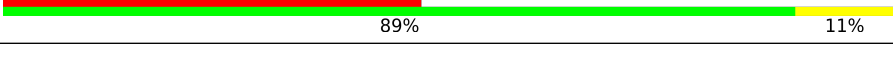
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	8855 (3.40 - 4.40)

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	471	20% 88% 10% .
1	E	471	19% 87% 10% .
1	I	471	20% 88% 10% .
2	B	153	46% 82% 7% 11%

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Mol	Chain	Length	Quality of chain
2	F	153	
2	J	153	
3	C	121	
3	G	121	
3	K	121	
4	D	101	
4	H	101	
4	L	101	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 38283 atoms, of which 18936 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 CH848 DE3 SOSIP gp120.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	461	Total	C	H	N	O	S	0	0
			7205	2273	3573	642	688	29		
1	E	461	Total	C	H	N	O	S	0	0
			7205	2273	3573	642	688	29		
1	I	461	Total	C	H	N	O	S	0	0
			7205	2273	3573	642	688	29		

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	-	expression tag	UNP A0A1W6IPB2
A	2	GLU	-	expression tag	UNP A0A1W6IPB2
A	3	ASN	-	expression tag	UNP A0A1W6IPB2
A	163	CYS	VAL	conflict	UNP A0A1W6IPB2
A	391	CYS	ALA	conflict	UNP A0A1W6IPB2
A	448	LYS	GLU	conflict	UNP A0A1W6IPB2
A	450	GLU	GLN	conflict	UNP A0A1W6IPB2
A	454	VAL	ILE	conflict	UNP A0A1W6IPB2
A	458	ARG	GLY	conflict	UNP A0A1W6IPB2
A	459	CYS	ALA	conflict	UNP A0A1W6IPB2
A	465	GLY	GLU	conflict	UNP A0A1W6IPB2
A	467	ARG	GLU	conflict	UNP A0A1W6IPB2
A	468	ARG	LYS	conflict	UNP A0A1W6IPB2
A	470	ARG	-	expression tag	UNP A0A1W6IPB2
A	471	ARG	-	expression tag	UNP A0A1W6IPB2
E	1	ALA	-	expression tag	UNP A0A1W6IPB2
E	2	GLU	-	expression tag	UNP A0A1W6IPB2
E	3	ASN	-	expression tag	UNP A0A1W6IPB2
E	163	CYS	VAL	conflict	UNP A0A1W6IPB2
E	391	CYS	ALA	conflict	UNP A0A1W6IPB2
E	448	LYS	GLU	conflict	UNP A0A1W6IPB2
E	450	GLU	GLN	conflict	UNP A0A1W6IPB2
E	454	VAL	ILE	conflict	UNP A0A1W6IPB2
E	458	ARG	GLY	conflict	UNP A0A1W6IPB2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	459	CYS	ALA	conflict	UNP A0A1W6IPB2
E	465	GLY	GLU	conflict	UNP A0A1W6IPB2
E	467	ARG	GLU	conflict	UNP A0A1W6IPB2
E	468	ARG	LYS	conflict	UNP A0A1W6IPB2
E	470	ARG	-	expression tag	UNP A0A1W6IPB2
E	471	ARG	-	expression tag	UNP A0A1W6IPB2
I	1	ALA	-	expression tag	UNP A0A1W6IPB2
I	2	GLU	-	expression tag	UNP A0A1W6IPB2
I	3	ASN	-	expression tag	UNP A0A1W6IPB2
I	163	CYS	VAL	conflict	UNP A0A1W6IPB2
I	391	CYS	ALA	conflict	UNP A0A1W6IPB2
I	448	LYS	GLU	conflict	UNP A0A1W6IPB2
I	450	GLU	GLN	conflict	UNP A0A1W6IPB2
I	454	VAL	ILE	conflict	UNP A0A1W6IPB2
I	458	ARG	GLY	conflict	UNP A0A1W6IPB2
I	459	CYS	ALA	conflict	UNP A0A1W6IPB2
I	465	GLY	GLU	conflict	UNP A0A1W6IPB2
I	467	ARG	GLU	conflict	UNP A0A1W6IPB2
I	468	ARG	LYS	conflict	UNP A0A1W6IPB2
I	470	ARG	-	expression tag	UNP A0A1W6IPB2
I	471	ARG	-	expression tag	UNP A0A1W6IPB2

- Molecule 2 is a protein called CH848 DE3 SOSIP gp41.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	136	Total	C	H	N	O	S	0	0
			2127	677	1060	183	201	6		
2	F	136	Total	C	H	N	O	S	0	0
			2127	677	1060	183	201	6		
2	J	136	Total	C	H	N	O	S	0	0
			2127	677	1060	183	201	6		

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	473	VAL	ALA	conflict	UNP A0A1W6IPB2
B	475	ILE	LEU	conflict	UNP A0A1W6IPB2
B	478	VAL	LEU	conflict	UNP A0A1W6IPB2
B	495	MET	ILE	conflict	UNP A0A1W6IPB2
B	503	ASN	GLN	conflict	UNP A0A1W6IPB2
B	527	LYS	GLN	conflict	UNP A0A1W6IPB2
B	543	VAL	LEU	conflict	UNP A0A1W6IPB2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	548	ARG	LYS	conflict	UNP A0A1W6IPB2
B	555	ILE	MET	conflict	UNP A0A1W6IPB2
B	565	CYS	THR	conflict	UNP A0A1W6IPB2
B	572	SER	THR	conflict	UNP A0A1W6IPB2
B	577	ARG	LYS	conflict	UNP A0A1W6IPB2
B	578	ASN	SER	conflict	UNP A0A1W6IPB2
B	579	LEU	GLU	conflict	UNP A0A1W6IPB2
B	580	SER	THR	conflict	UNP A0A1W6IPB2
B	581	GLU	ASP	conflict	UNP A0A1W6IPB2
B	589	LEU	MET	conflict	UNP A0A1W6IPB2
B	592	ASP	GLU	conflict	UNP A0A1W6IPB2
B	593	LYS	ARG	conflict	UNP A0A1W6IPB2
B	600	GLN	GLU	conflict	UNP A0A1W6IPB2
B	601	ILE	THR	conflict	UNP A0A1W6IPB2
B	604	GLY	LYS	conflict	UNP A0A1W6IPB2
B	608	GLU	ASP	conflict	UNP A0A1W6IPB2
B	615	LYS	ARG	conflict	UNP A0A1W6IPB2
F	473	VAL	ALA	conflict	UNP A0A1W6IPB2
F	475	ILE	LEU	conflict	UNP A0A1W6IPB2
F	478	VAL	LEU	conflict	UNP A0A1W6IPB2
F	495	MET	ILE	conflict	UNP A0A1W6IPB2
F	503	ASN	GLN	conflict	UNP A0A1W6IPB2
F	527	LYS	GLN	conflict	UNP A0A1W6IPB2
F	543	VAL	LEU	conflict	UNP A0A1W6IPB2
F	548	ARG	LYS	conflict	UNP A0A1W6IPB2
F	555	ILE	MET	conflict	UNP A0A1W6IPB2
F	565	CYS	THR	conflict	UNP A0A1W6IPB2
F	572	SER	THR	conflict	UNP A0A1W6IPB2
F	577	ARG	LYS	conflict	UNP A0A1W6IPB2
F	578	ASN	SER	conflict	UNP A0A1W6IPB2
F	579	LEU	GLU	conflict	UNP A0A1W6IPB2
F	580	SER	THR	conflict	UNP A0A1W6IPB2
F	581	GLU	ASP	conflict	UNP A0A1W6IPB2
F	589	LEU	MET	conflict	UNP A0A1W6IPB2
F	592	ASP	GLU	conflict	UNP A0A1W6IPB2
F	593	LYS	ARG	conflict	UNP A0A1W6IPB2
F	600	GLN	GLU	conflict	UNP A0A1W6IPB2
F	601	ILE	THR	conflict	UNP A0A1W6IPB2
F	604	GLY	LYS	conflict	UNP A0A1W6IPB2
F	608	GLU	ASP	conflict	UNP A0A1W6IPB2
F	615	LYS	ARG	conflict	UNP A0A1W6IPB2
J	473	VAL	ALA	conflict	UNP A0A1W6IPB2

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Chain	Residue	Modelled	Actual	Comment	Reference
J	475	ILE	LEU	conflict	UNP A0A1W6IPB2
J	478	VAL	LEU	conflict	UNP A0A1W6IPB2
J	495	MET	ILE	conflict	UNP A0A1W6IPB2
J	503	ASN	GLN	conflict	UNP A0A1W6IPB2
J	527	LYS	GLN	conflict	UNP A0A1W6IPB2
J	543	VAL	LEU	conflict	UNP A0A1W6IPB2
J	548	ARG	LYS	conflict	UNP A0A1W6IPB2
J	555	ILE	MET	conflict	UNP A0A1W6IPB2
J	565	CYS	THR	conflict	UNP A0A1W6IPB2
J	572	SER	THR	conflict	UNP A0A1W6IPB2
J	577	ARG	LYS	conflict	UNP A0A1W6IPB2
J	578	ASN	SER	conflict	UNP A0A1W6IPB2
J	579	LEU	GLU	conflict	UNP A0A1W6IPB2
J	580	SER	THR	conflict	UNP A0A1W6IPB2
J	581	GLU	ASP	conflict	UNP A0A1W6IPB2
J	589	LEU	MET	conflict	UNP A0A1W6IPB2
J	592	ASP	GLU	conflict	UNP A0A1W6IPB2
J	593	LYS	ARG	conflict	UNP A0A1W6IPB2
J	600	GLN	GLU	conflict	UNP A0A1W6IPB2
J	601	ILE	THR	conflict	UNP A0A1W6IPB2
J	604	GLY	LYS	conflict	UNP A0A1W6IPB2
J	608	GLU	ASP	conflict	UNP A0A1W6IPB2
J	615	LYS	ARG	conflict	UNP A0A1W6IPB2

- Molecule 3 is a protein called VRC01 Fab Heavy Chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	121	Total	C	H	N	O	S	0	0
			1904	609	937	175	175	8		
3	G	121	Total	C	H	N	O	S	0	0
			1904	609	937	175	175	8		
3	K	121	Total	C	H	N	O	S	0	0
			1904	609	937	175	175	8		

- Molecule 4 is a protein called VRC01 Fab Light Chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	101	Total	C	H	N	O	S	0	0
			1525	494	742	133	154	2		
4	H	101	Total	C	H	N	O	S	0	0
			1525	494	742	133	154	2		

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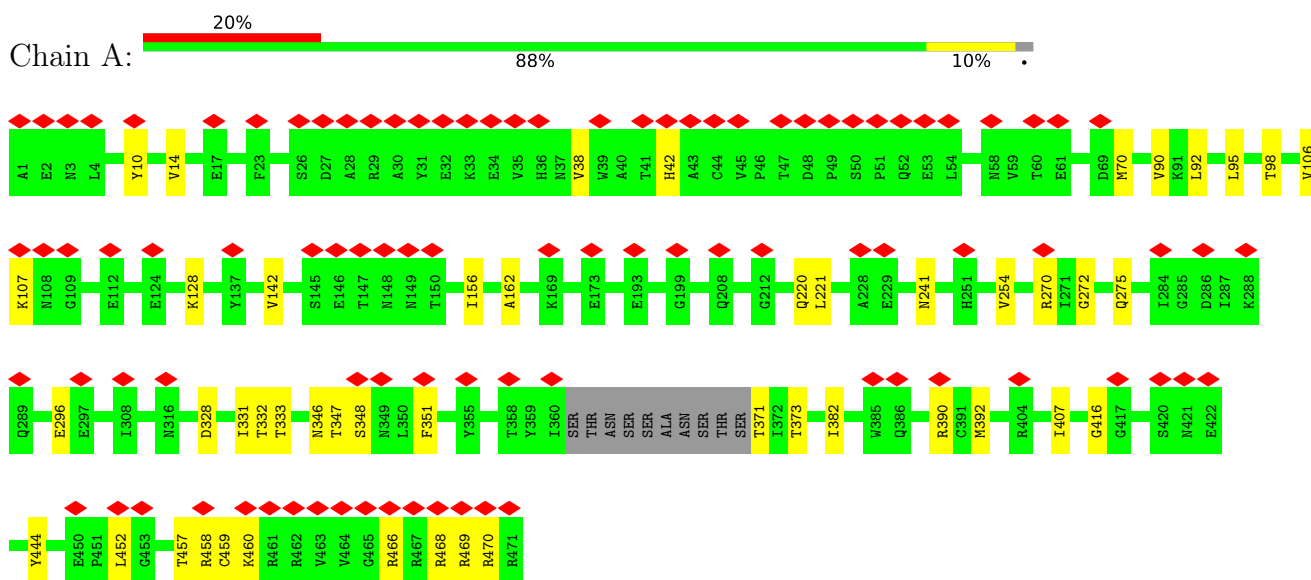
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Mol	Chain	Residues	Atoms						AltConf	Trace
4	L	101	Total	C	H	N	O	S	0	0
			1525	494	742	133	154	2		

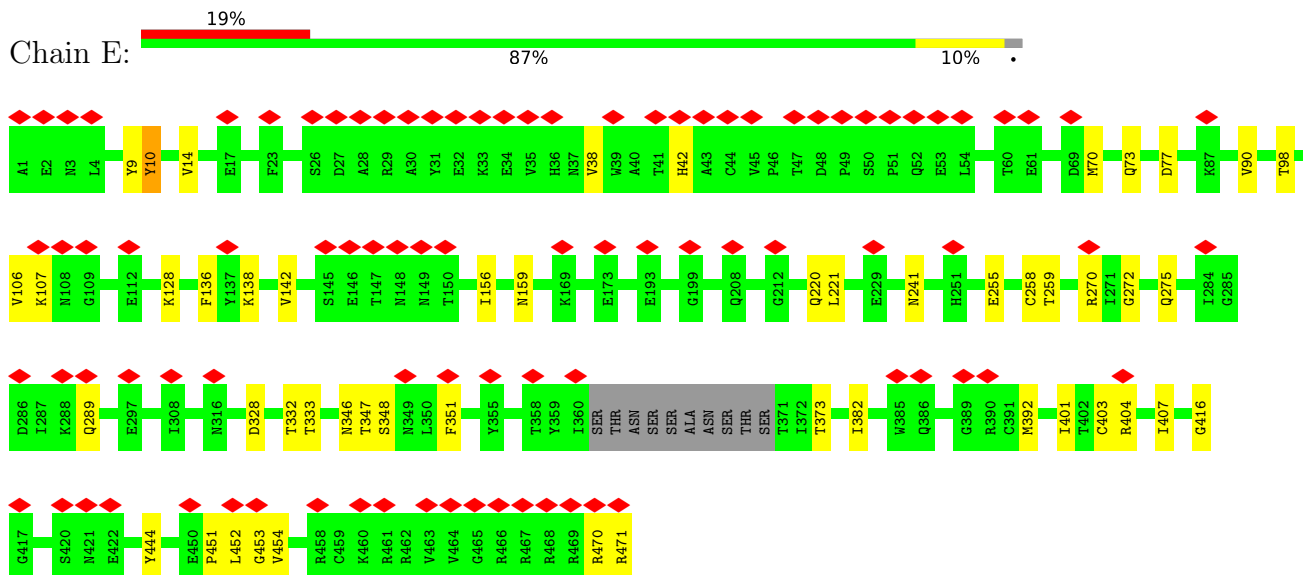
3 Residue-property plots [i](#)

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.

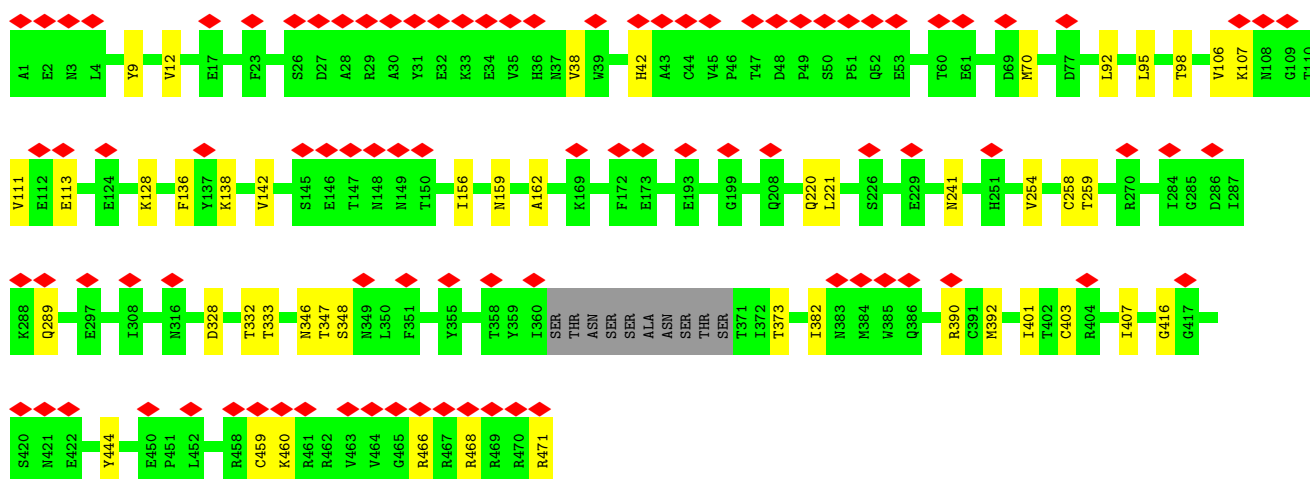
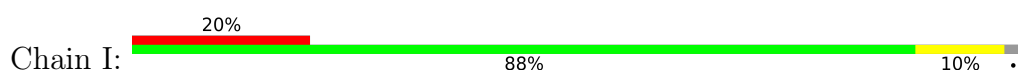
• Molecule 1: CH848 DE3 SOSIP gp120



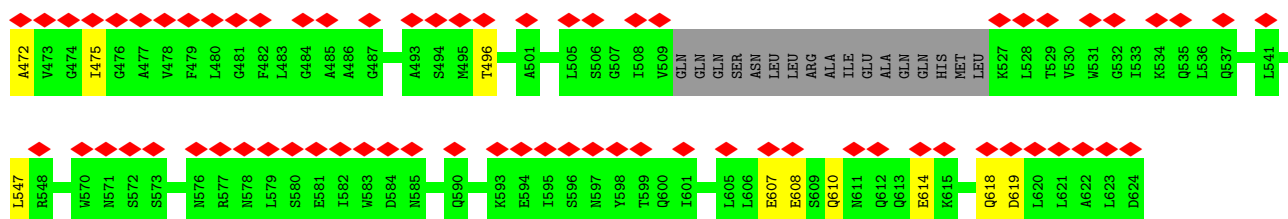
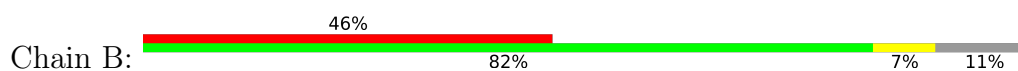
• Molecule 1: CH848 DE3 SOSIP gp120



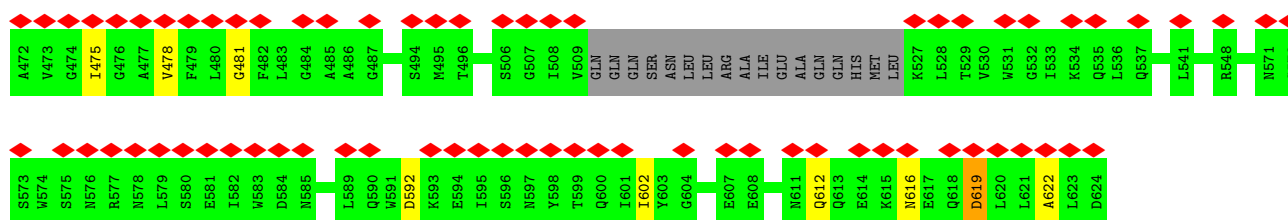
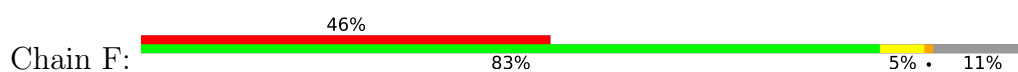
• Molecule 1: CH848 DE3 SOSIP gp120



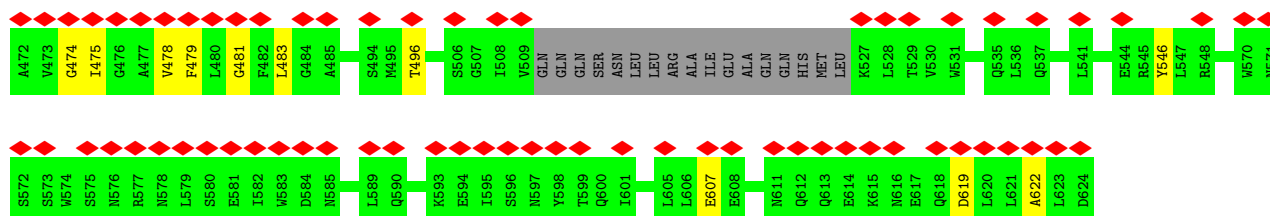
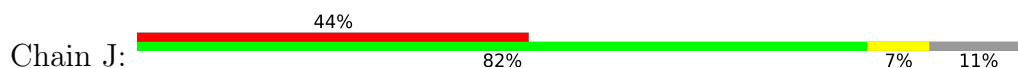
• Molecule 2: CH848 DE3 SOSIP gp41



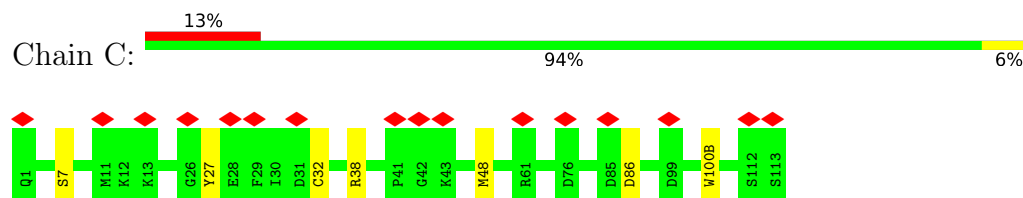
• Molecule 2: CH848 DE3 SOSIP gp41



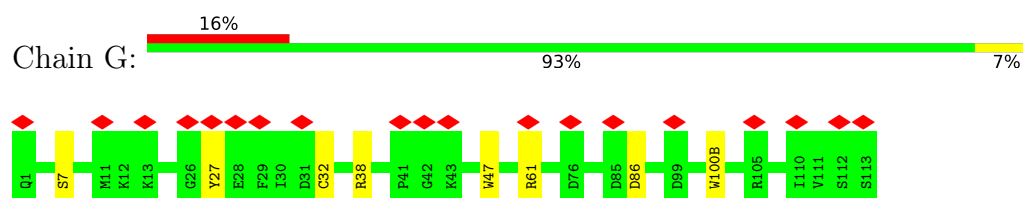
• Molecule 2: CH848 DE3 SOSIP gp41



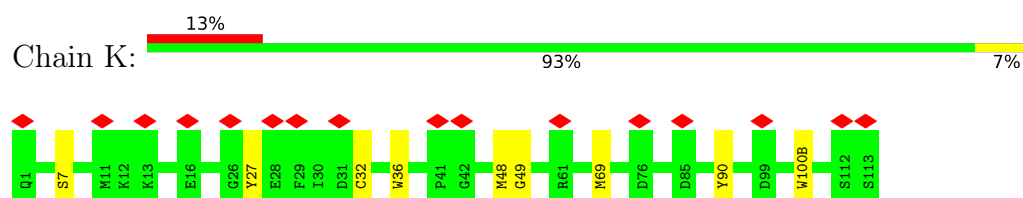
- Molecule 3: VRC01 Fab Heavy Chain



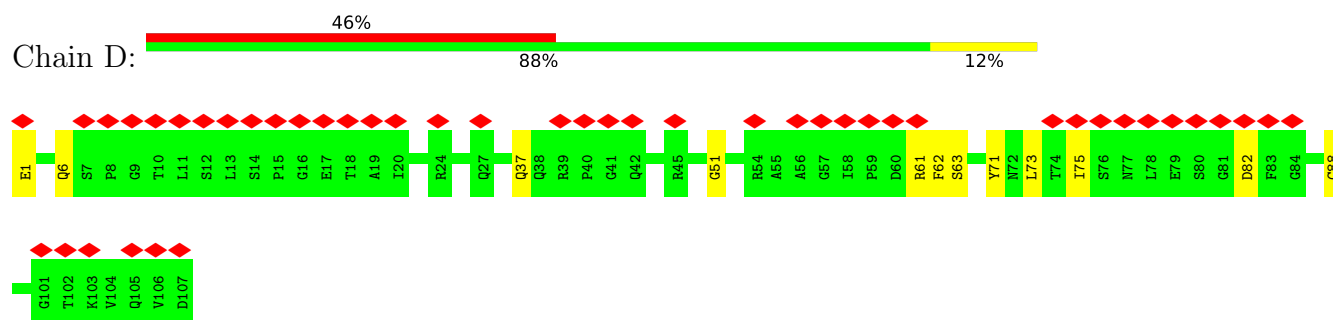
- Molecule 3: VRC01 Fab Heavy Chain



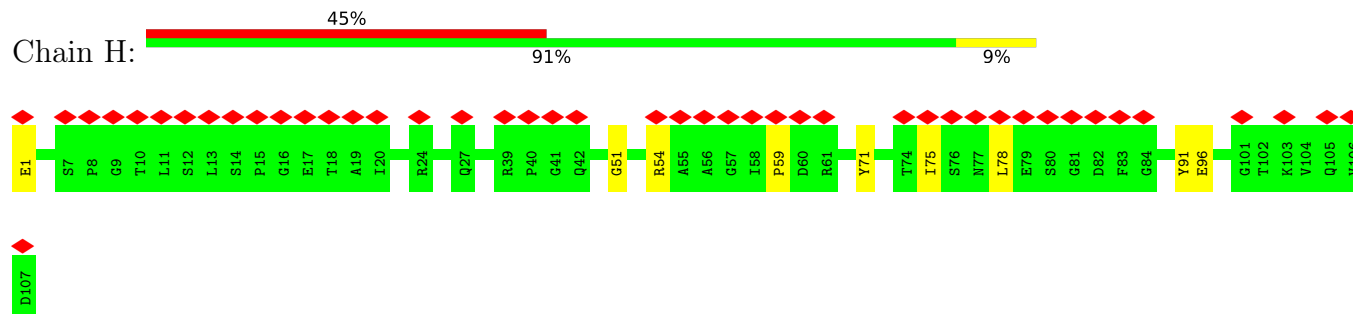
- Molecule 3: VRC01 Fab Heavy Chain



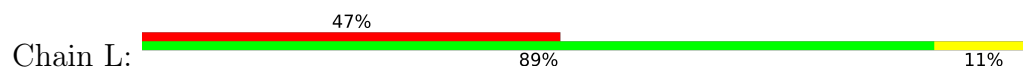
- Molecule 4: VRC01 Fab Light Chain

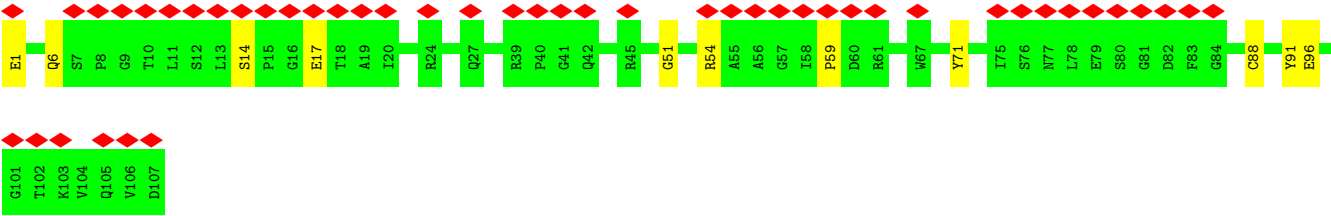


- Molecule 4: VRC01 Fab Light Chain



- Molecule 4: VRC01 Fab Light Chain





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	220881	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	420	Depositor
Maximum defocus (nm)	4710	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	13.170	Depositor
Minimum map value	-8.054	Depositor
Average map value	0.008	Depositor
Map value standard deviation	0.261	Depositor
Recommended contour level	2.1	Depositor
Map size ($\text{\AA}$)	324.0, 324.0, 324.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.08, 1.08, 1.08	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	0.20	0/3708	0.49	4/5036 (0.1%)
1	E	0.20	0/3708	0.51	4/5036 (0.1%)
1	I	0.20	0/3708	0.50	4/5036 (0.1%)
2	B	0.17	0/1085	0.44	0/1470
2	F	0.18	0/1085	0.43	0/1470
2	J	0.17	0/1085	0.42	0/1470
3	C	0.20	0/992	0.50	2/1343 (0.1%)
3	G	0.19	0/992	0.49	2/1343 (0.1%)
3	K	0.19	0/992	0.48	2/1343 (0.1%)
4	D	0.18	0/802	0.37	0/1089
4	H	0.18	0/802	0.41	0/1089
4	L	0.18	0/802	0.37	0/1089
All	All	0.19	0/19761	0.47	18/26814 (0.1%)

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
1	A	0	3
1	E	0	5
1	I	0	1
2	F	0	1
All	All	0	10

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	HIS	CA-C-N	6.00	132.50	121.70
1	A	42	HIS	C-N-CA	6.00	132.50	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	42	HIS	CA-C-N	5.96	132.43	121.70
1	E	42	HIS	C-N-CA	5.96	132.43	121.70
1	I	42	HIS	CA-C-N	5.80	132.14	121.70
1	I	42	HIS	C-N-CA	5.80	132.14	121.70
3	C	7	SER	CA-C-N	5.34	131.31	121.70
3	C	7	SER	C-N-CA	5.34	131.31	121.70
3	G	7	SER	CA-C-N	5.33	131.29	121.70
3	G	7	SER	C-N-CA	5.33	131.29	121.70
1	I	416	GLY	CA-C-N	5.31	131.26	121.70
1	I	416	GLY	C-N-CA	5.31	131.26	121.70
1	E	416	GLY	CA-C-N	5.14	130.96	121.70
1	E	416	GLY	C-N-CA	5.14	130.96	121.70
3	K	7	SER	CA-C-N	5.11	130.90	121.70
3	K	7	SER	C-N-CA	5.11	130.90	121.70
1	A	416	GLY	CA-C-N	5.08	130.84	121.70
1	A	416	GLY	C-N-CA	5.08	130.84	121.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	TYR	Peptide
1	A	351	PHE	Peptide
1	A	38	VAL	Peptide
1	E	10	TYR	Peptide
1	E	351	PHE	Peptide
1	E	38	VAL	Peptide
1	E	451	PRO	Peptide
1	E	9	TYR	Peptide
2	F	619	ASP	Peptide
1	I	38	VAL	Peptide

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	3632	3573	3577	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3632	3573	3577	29	0
1	I	3632	3573	3577	27	0
2	B	1067	1060	1058	8	0
2	F	1067	1060	1058	6	0
2	J	1067	1060	1058	9	0
3	C	967	937	939	4	0
3	G	967	937	939	4	0
3	K	967	937	939	5	0
4	D	783	742	743	7	0
4	H	783	742	743	5	0
4	L	783	742	743	6	0
All	All	19347	18936	18951	119	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 3.

All (119) 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:A:382:ILE:HD11	1:A:392:MET:HE2	1.69	0.74
4:L:91:TYR:O	4:L:96:GLU:N	2.22	0.73
1:A:162:ALA:O	1:A:390:ARG:NH2	2.22	0.73
1:E:382:ILE:HD11	1:E:392:MET:HE2	1.71	0.72
1:I:162:ALA:O	1:I:390:ARG:NH2	2.23	0.71
1:E:73:GLN:NE2	1:E:77:ASP:OD2	2.25	0.69
1:E:259:THR:OG1	1:E:401:ILE:O	2.05	0.69
1:I:382:ILE:HD11	1:I:392:MET:HE2	1.75	0.69
4:H:54:ARG:NH2	4:H:59:PRO:O	2.27	0.68
1:I:259:THR:OG1	1:I:401:ILE:O	2.04	0.68
1:I:92:LEU:HD12	1:I:95:LEU:HD12	1.74	0.68
4:L:51:GLY:O	4:L:71:TYR:OH	2.10	0.68
4:H:51:GLY:O	4:H:71:TYR:OH	2.10	0.67
4:H:91:TYR:O	4:H:96:GLU:N	2.28	0.66
1:A:98:THR:OG1	1:I:128:LYS:NZ	2.29	0.66
1:A:128:LYS:NZ	1:E:98:THR:OG1	2.29	0.66
4:D:37:GLN:NE2	4:D:82:ASP:O	2.30	0.65
3:G:47:TRP:O	3:G:61:ARG:NH1	2.32	0.63
1:A:92:LEU:HD12	1:A:95:LEU:HD12	1.80	0.62
1:A:270:ARG:NH1	1:A:272:GLY:O	2.33	0.62
3:C:27:TYR:OH	3:C:32:CYS:SG	2.59	0.61
1:A:14:VAL:HG21	1:A:452:LEU:HD11	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:14:VAL:HG21	1:E:452:LEU:HD11	1.83	0.60
1:A:296:GLU:OE2	1:A:371:THR:N	2.36	0.59
2:F:475:ILE:HG21	2:J:607:GLU:HB3	1.83	0.59
1:A:142:VAL:HG23	1:A:156:ILE:HG22	1.84	0.58
4:L:54:ARG:NH2	4:L:59:PRO:O	2.36	0.58
1:E:258:CYS:O	1:E:403:CYS:N	2.38	0.57
4:D:6:GLN:OE1	4:D:88:CYS:N	2.38	0.56
1:E:221:LEU:HD22	1:E:407:ILE:HD11	1.87	0.56
1:E:454:VAL:HG21	2:F:602:ILE:HG21	1.87	0.56
1:A:221:LEU:HD22	1:A:407:ILE:HD11	1.88	0.55
3:K:48:MET:HE1	3:K:90:TYR:CD2	2.41	0.55
3:G:27:TYR:OH	3:G:32:CYS:SG	2.65	0.55
1:E:10:TYR:OH	1:E:453:GLY:O	2.19	0.55
1:E:128:LYS:NZ	1:I:98:THR:OG1	2.40	0.55
1:I:258:CYS:O	1:I:403:CYS:N	2.38	0.55
1:A:470:ARG:O	2:F:622:ALA:N	2.36	0.54
2:B:475:ILE:HA	2:B:496:THR:HG23	1.90	0.53
1:I:9:TYR:O	1:I:12:VAL:HG23	2.08	0.53
1:I:466:ARG:O	1:I:468:ARG:NH1	2.42	0.53
1:E:270:ARG:NH1	1:E:272:GLY:O	2.42	0.53
1:I:346:ASN:O	1:I:347:THR:OG1	2.24	0.52
1:I:142:VAL:HG23	1:I:156:ILE:HG22	1.90	0.52
2:B:619:ASP:O	1:I:471:ARG:NH2	2.43	0.52
1:A:348:SER:OG	1:A:373:THR:O	2.23	0.52
3:K:27:TYR:OH	3:K:32:CYS:SG	2.68	0.51
1:E:348:SER:OG	1:E:373:THR:O	2.24	0.51
4:L:14:SER:N	4:L:17:GLU:OE2	2.43	0.51
2:B:607:GLU:HB3	2:J:475:ILE:HD12	1.92	0.51
1:I:254:VAL:O	1:I:407:ILE:HG22	2.10	0.51
1:E:142:VAL:HG23	1:E:156:ILE:HG22	1.91	0.51
2:F:612:GLN:O	2:F:616:ASN:ND2	2.43	0.51
1:A:92:LEU:HD12	1:A:95:LEU:CD1	2.41	0.50
1:E:241:ASN:OD1	3:G:100(B):TRP:NE1	2.45	0.49
1:I:220:GLN:O	1:I:221:LEU:HD23	2.13	0.48
1:I:332:THR:HG23	1:I:333:THR:HG23	1.95	0.48
1:A:466:ARG:O	1:A:468:ARG:NH1	2.46	0.48
1:E:136:PHE:O	1:E:138:LYS:N	2.45	0.48
2:B:547:LEU:HD12	2:J:546:TYR:CE2	2.49	0.48
1:I:92:LEU:HD12	1:I:95:LEU:CD1	2.41	0.48
4:D:62:PHE:HD1	4:D:75:ILE:HG22	1.78	0.47
1:I:221:LEU:HD22	1:I:407:ILE:HD11	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:THR:HG23	1:A:333:THR:HG23	1.96	0.47
1:E:346:ASN:O	1:E:347:THR:OG1	2.24	0.47
1:A:241:ASN:OD1	3:C:100(B):TRP:NE1	2.47	0.47
1:A:469:ARG:HG2	1:A:470:ARG:H	1.79	0.47
1:A:270:ARG:NH2	1:E:159:ASN:O	2.49	0.46
1:I:348:SER:OG	1:I:373:THR:O	2.22	0.46
1:E:220:GLN:O	1:E:221:LEU:HD23	2.16	0.46
4:D:51:GLY:O	4:D:71:TYR:OH	2.25	0.46
4:H:75:ILE:HD12	4:H:78:LEU:HD13	1.97	0.46
1:E:332:THR:HG23	1:E:333:THR:HG23	1.97	0.45
2:B:618:GLN:CA	1:I:471:ARG:HA	2.47	0.45
1:A:346:ASN:O	1:A:347:THR:OG1	2.23	0.45
2:B:610:GLN:O	2:B:614:GLU:N	2.47	0.45
2:J:478:VAL:O	2:J:481:GLY:N	2.50	0.45
1:A:220:GLN:O	1:A:221:LEU:HD23	2.17	0.45
4:D:61:ARG:NH1	4:D:82:ASP:OD2	2.50	0.45
1:E:470:ARG:O	2:J:622:ALA:N	2.42	0.45
4:L:1:GLU:OE1	4:L:1:GLU:N	2.50	0.45
2:B:608:GLU:HA	2:J:475:ILE:HD11	1.99	0.44
4:H:1:GLU:N	4:H:1:GLU:OE1	2.50	0.44
1:I:70:MET:HE1	1:I:444:TYR:CB	2.48	0.44
1:A:90:VAL:HG23	1:A:275:GLN:HG3	2.00	0.44
3:C:38:ARG:HB2	3:C:48:MET:HE3	1.99	0.43
1:A:220:GLN:NE2	1:A:331:ILE:O	2.50	0.43
1:I:459:CYS:SG	1:I:460:LYS:N	2.92	0.43
1:A:70:MET:HE1	1:A:444:TYR:CB	2.48	0.43
3:K:36:TRP:CE3	3:K:69:MET:HE3	2.54	0.43
4:D:1:GLU:OE1	4:D:1:GLU:N	2.50	0.43
1:A:254:VAL:O	1:A:407:ILE:HG22	2.17	0.43
4:D:63:SER:O	4:D:73:LEU:HD12	2.18	0.43
1:E:270:ARG:NH2	1:I:159:ASN:O	2.52	0.43
1:I:136:PHE:O	1:I:138:LYS:N	2.47	0.43
1:I:241:ASN:OD1	3:K:100(B):TRP:NE1	2.50	0.43
1:E:452:LEU:HD13	2:F:592:ASP:HB2	2.00	0.43
1:A:459:CYS:SG	1:A:460:LYS:N	2.93	0.42
3:C:38:ARG:NH2	3:C:86:ASP:OD1	2.52	0.42
1:E:106:VAL:HG22	1:E:107:LYS:H	1.83	0.42
2:B:472:ALA:O	2:B:496:THR:OG1	2.33	0.42
1:E:70:MET:HE1	1:E:444:TYR:CB	2.49	0.42
3:G:38:ARG:NH2	3:G:86:ASP:OD1	2.53	0.42
1:I:111:VAL:HG22	1:I:113:GLU:H	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:90:VAL:HG23	1:E:275:GLN:HG3	2.01	0.42
4:L:6:GLN:OE1	4:L:88:CYS:N	2.53	0.42
1:A:328:ASP:O	1:A:332:THR:HG22	2.20	0.42
1:I:106:VAL:HG22	1:I:107:LYS:H	1.84	0.42
1:E:255:GLU:OE2	1:E:404:ARG:NH1	2.50	0.42
1:E:328:ASP:O	1:E:332:THR:HG22	2.20	0.41
1:A:106:VAL:HG22	1:A:107:LYS:H	1.85	0.41
1:A:457:THR:HG22	1:A:458:ARG:H	1.84	0.41
1:I:328:ASP:O	1:I:332:THR:HG22	2.20	0.41
2:F:478:VAL:O	2:F:481:GLY:N	2.53	0.41
1:E:14:VAL:CG2	1:E:452:LEU:HD11	2.50	0.41
2:J:474:GLY:O	2:J:496:THR:OG1	2.21	0.40
2:J:479:PHE:CZ	2:J:483:LEU:HD11	2.56	0.40
1:E:471:ARG:NH2	2:J:619:ASP:O	2.54	0.40
3:K:49:GLY:O	3:K:69:MET:HE1	2.22	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	457/471 (97%)	425 (93%)	32 (7%)	0	100	100
1	E	457/471 (97%)	426 (93%)	30 (7%)	1 (0%)	43	74
1	I	457/471 (97%)	422 (92%)	34 (7%)	1 (0%)	43	74
2	B	132/153 (86%)	119 (90%)	13 (10%)	0	100	100
2	F	132/153 (86%)	120 (91%)	11 (8%)	1 (1%)	16	50
2	J	132/153 (86%)	119 (90%)	13 (10%)	0	100	100
3	C	119/121 (98%)	111 (93%)	8 (7%)	0	100	100
3	G	119/121 (98%)	111 (93%)	8 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	K	119/121 (98%)	110 (92%)	9 (8%)	0	100	100
4	D	97/101 (96%)	96 (99%)	1 (1%)	0	100	100
4	H	97/101 (96%)	95 (98%)	2 (2%)	0	100	100
4	L	97/101 (96%)	96 (99%)	1 (1%)	0	100	100
All	All	2415/2538 (95%)	2250 (93%)	162 (7%)	3 (0%)	49	80

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	619	ASP
1	E	289	GLN
1	I	289	GLN

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	409/418 (98%)	409 (100%)	0	100	100
1	E	409/418 (98%)	409 (100%)	0	100	100
1	I	409/418 (98%)	409 (100%)	0	100	100
2	B	114/129 (88%)	114 (100%)	0	100	100
2	F	114/129 (88%)	114 (100%)	0	100	100
2	J	114/129 (88%)	114 (100%)	0	100	100
3	C	103/103 (100%)	103 (100%)	0	100	100
3	G	103/103 (100%)	103 (100%)	0	100	100
3	K	103/103 (100%)	103 (100%)	0	100	100
4	D	84/84 (100%)	84 (100%)	0	100	100
4	H	84/84 (100%)	84 (100%)	0	100	100
4	L	84/84 (100%)	84 (100%)	0	100	100
All	All	2130/2202 (97%)	2130 (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 (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	GLN
1	A	289	GLN
1	A	346	ASN
1	A	400	ASN
3	C	58	ASN
1	E	73	GLN
1	E	346	ASN
2	F	503	ASN
1	I	73	GLN
1	I	346	ASN
2	J	567	ASN
2	J	578	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 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 ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	D	1
4	H	1
4	L	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	91:TYR	C	96:GLU	N	3.11
1	H	91:TYR	C	96:GLU	N	3.10
1	L	91:TYR	C	96:GLU	N	3.06

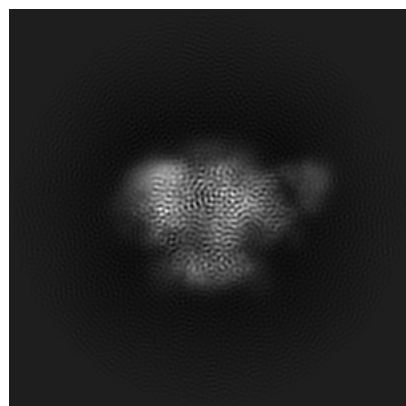
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-43228. These allow visual inspection of the internal detail of the map and identification of artifacts.

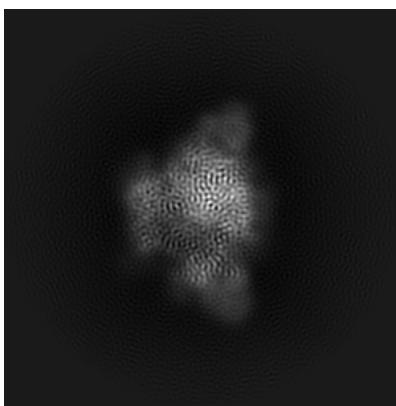
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

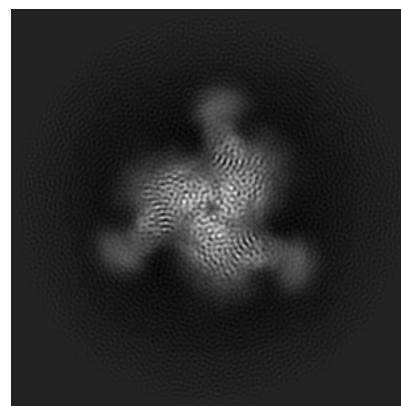
6.1.1 Primary map



X

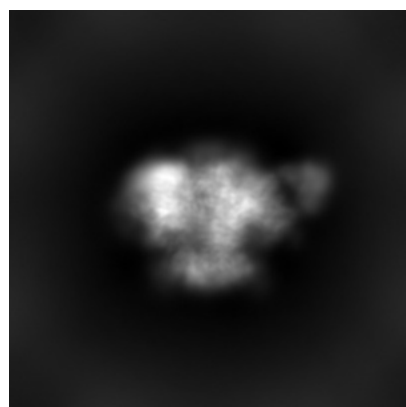


Y

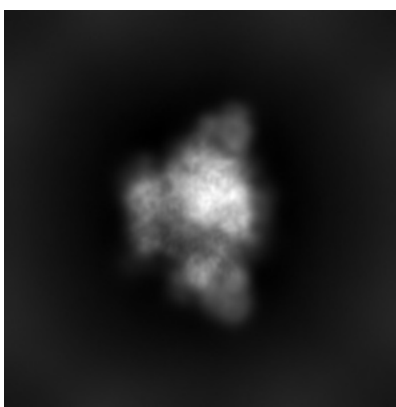


Z

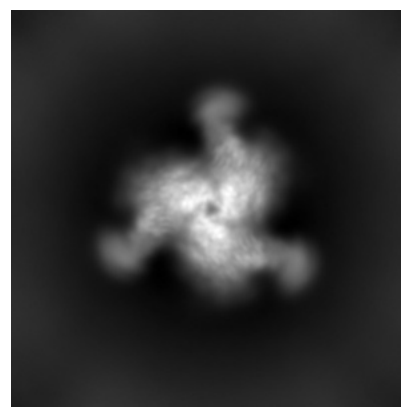
6.1.2 Raw map



X



Y

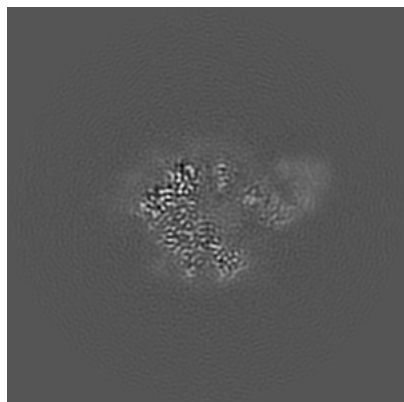


Z

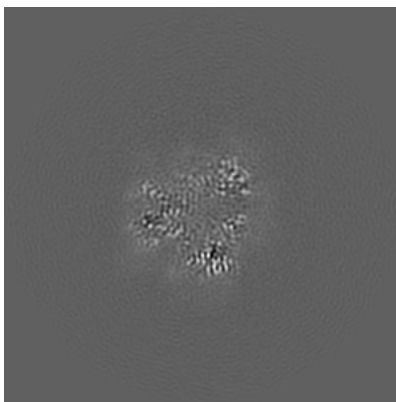
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

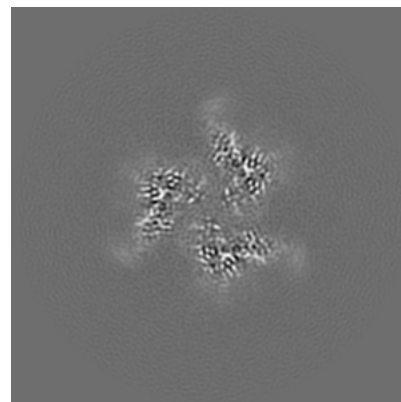
6.2.1 Primary map



X Index: 150

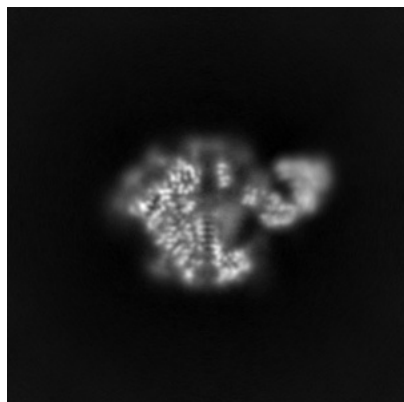


Y Index: 150

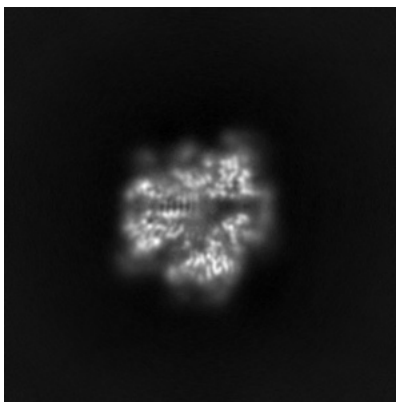


Z Index: 150

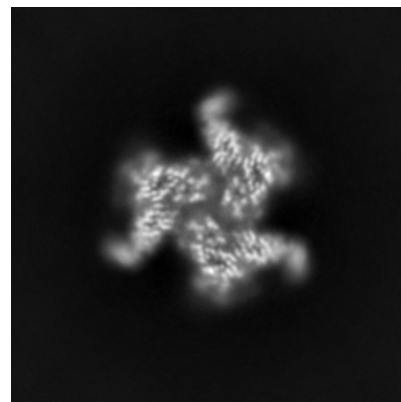
6.2.2 Raw map



X Index: 150



Y Index: 150

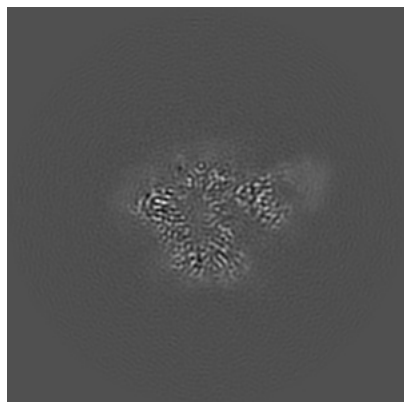


Z Index: 150

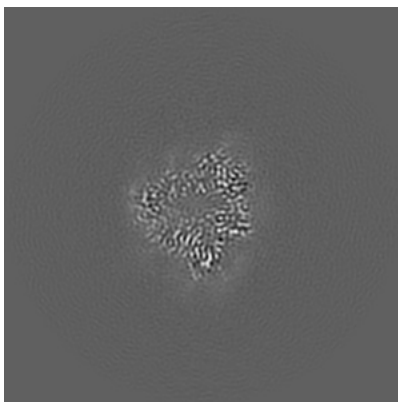
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

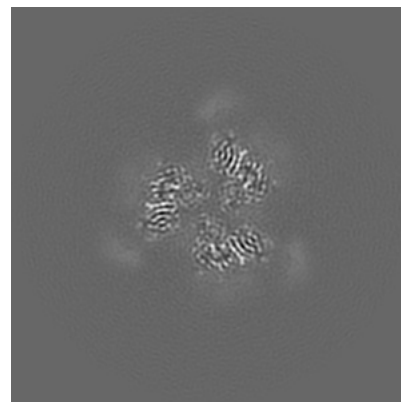
6.3.1 Primary map



X Index: 159

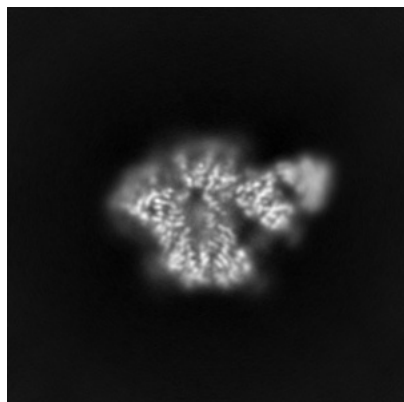


Y Index: 160

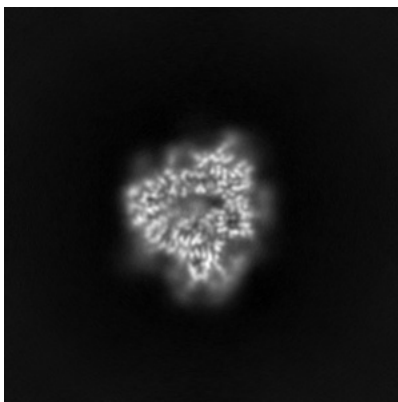


Z Index: 157

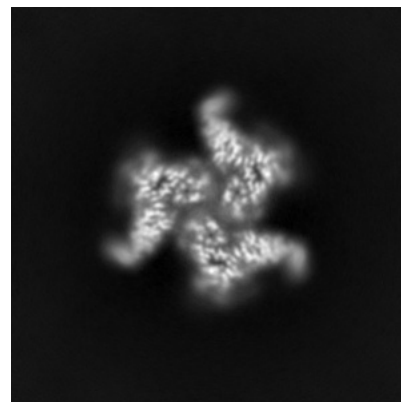
6.3.2 Raw map



X Index: 159



Y Index: 160

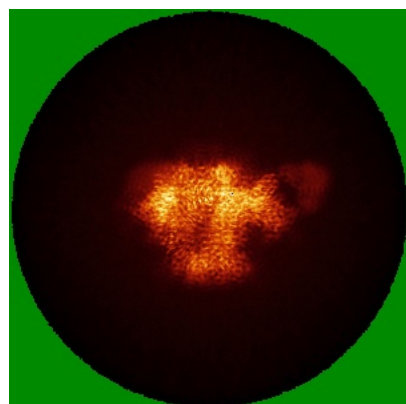


Z Index: 149

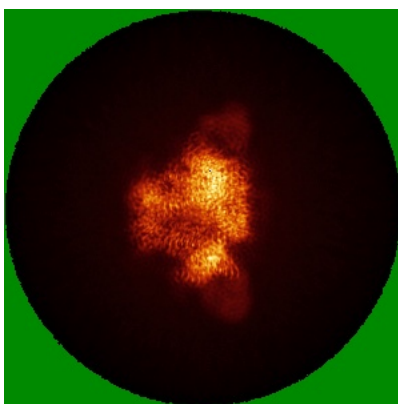
The images above show the largest variance slices of the map in three orthogonal directions.

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

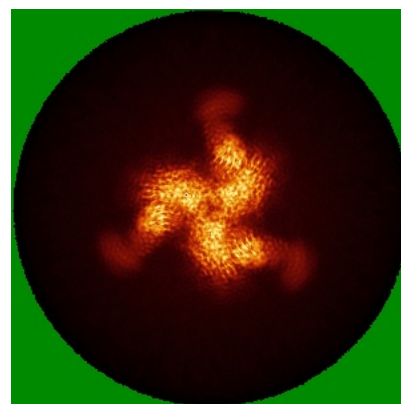
6.4.1 Primary map



X

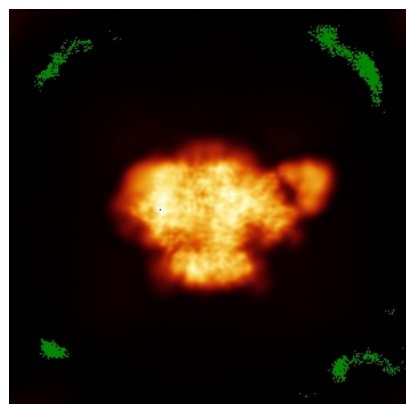


Y

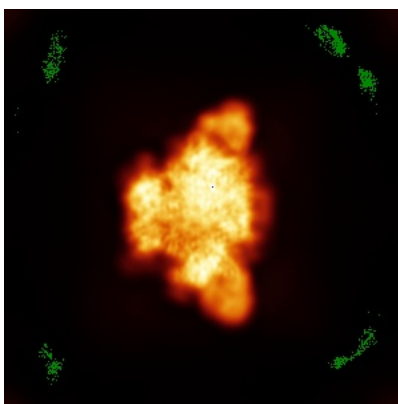


Z

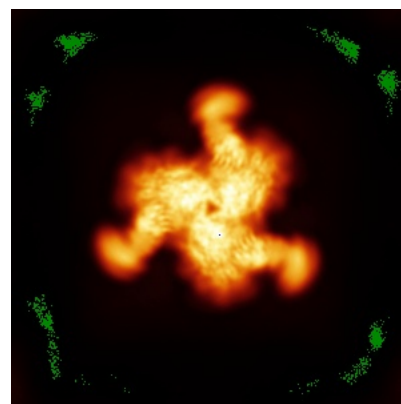
6.4.2 Raw map



X



Y

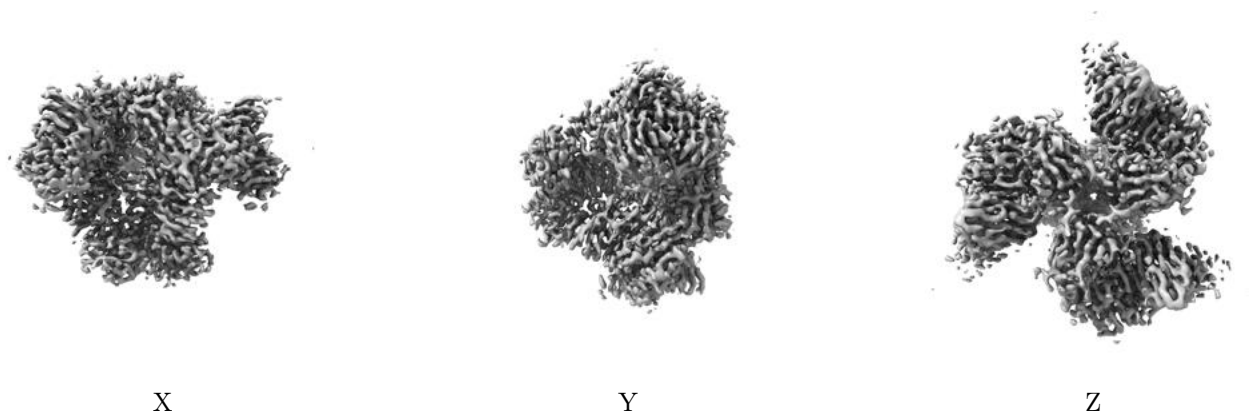


Z

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

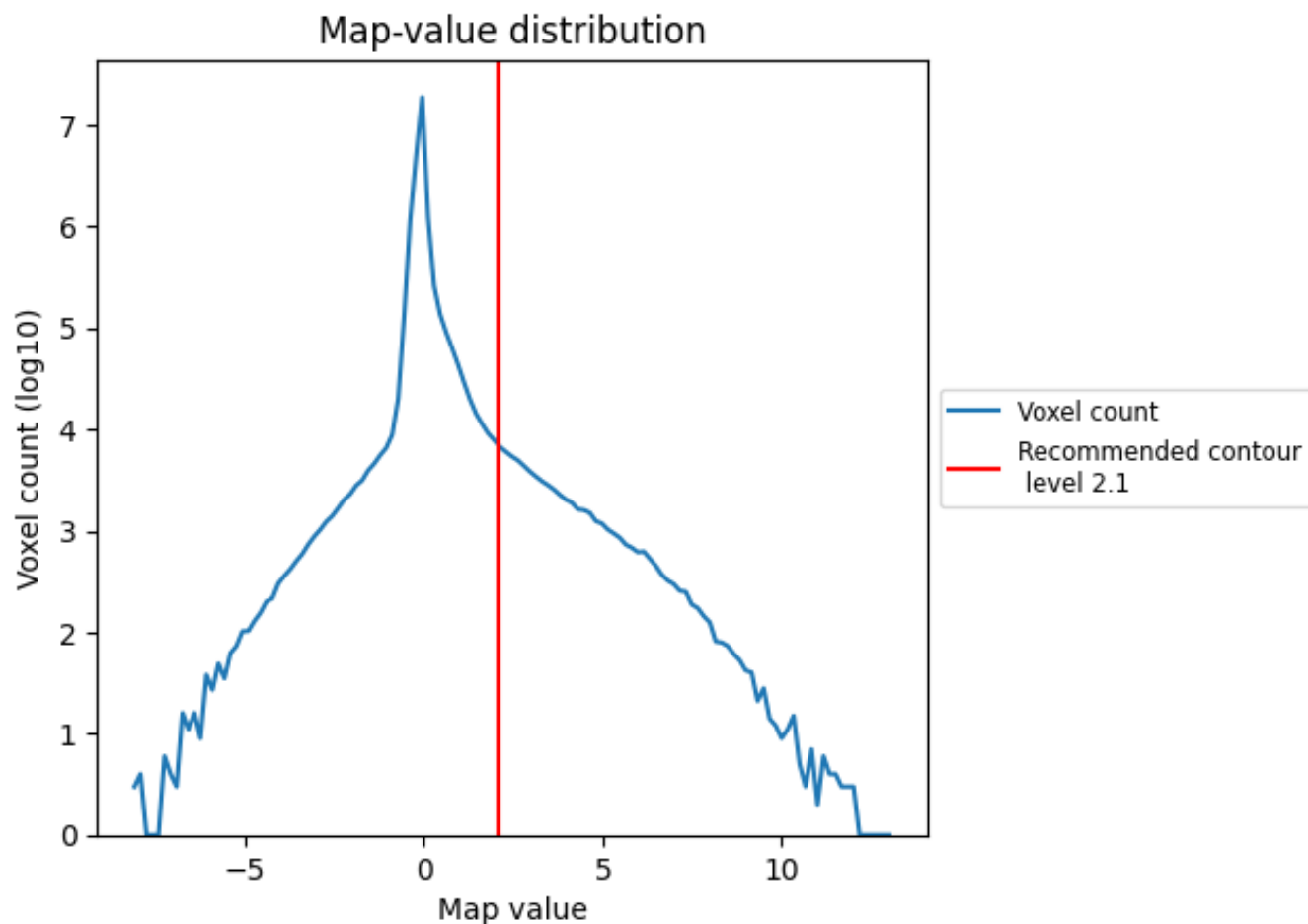
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

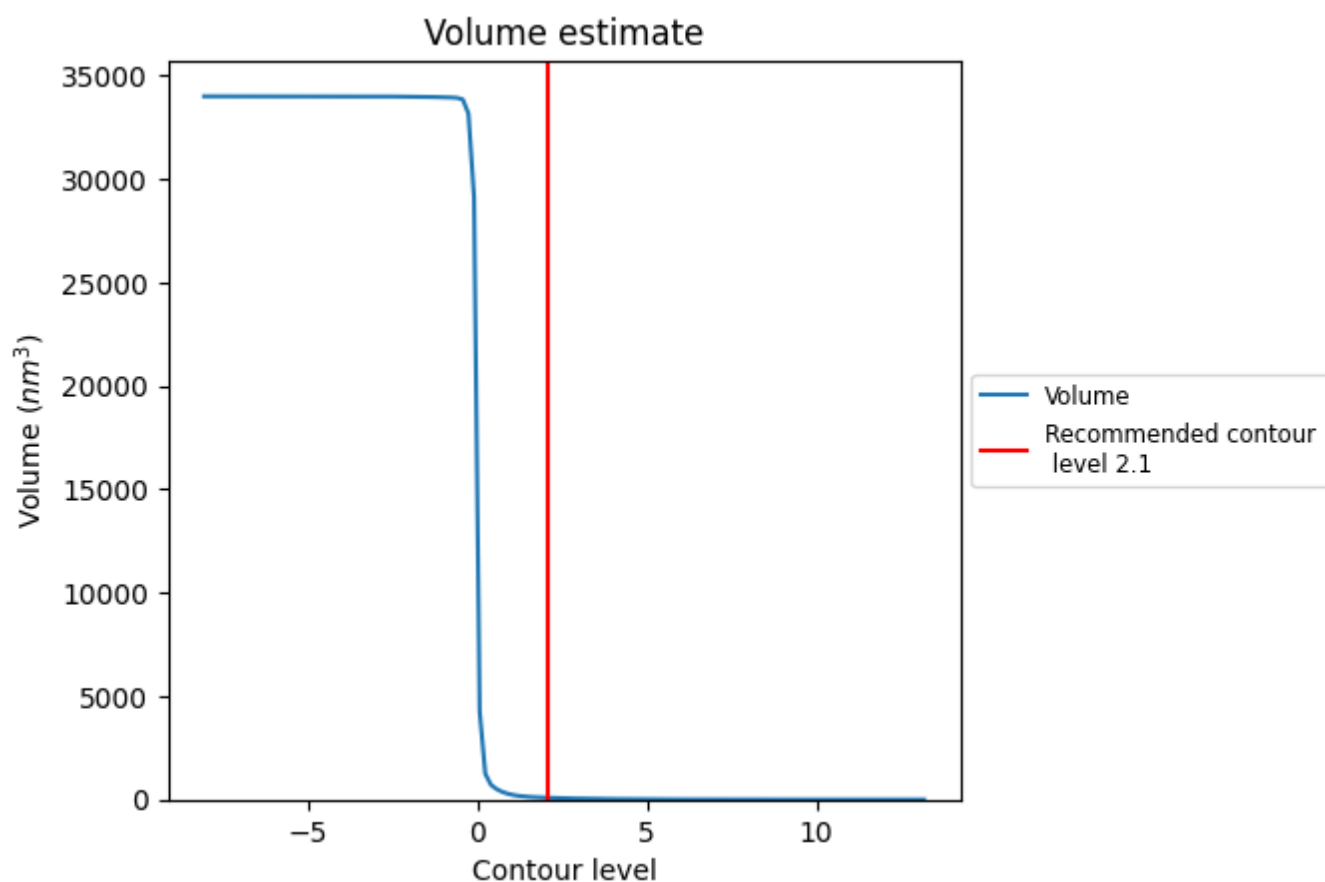
This section contains the results of statistical analysis of the map.

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.

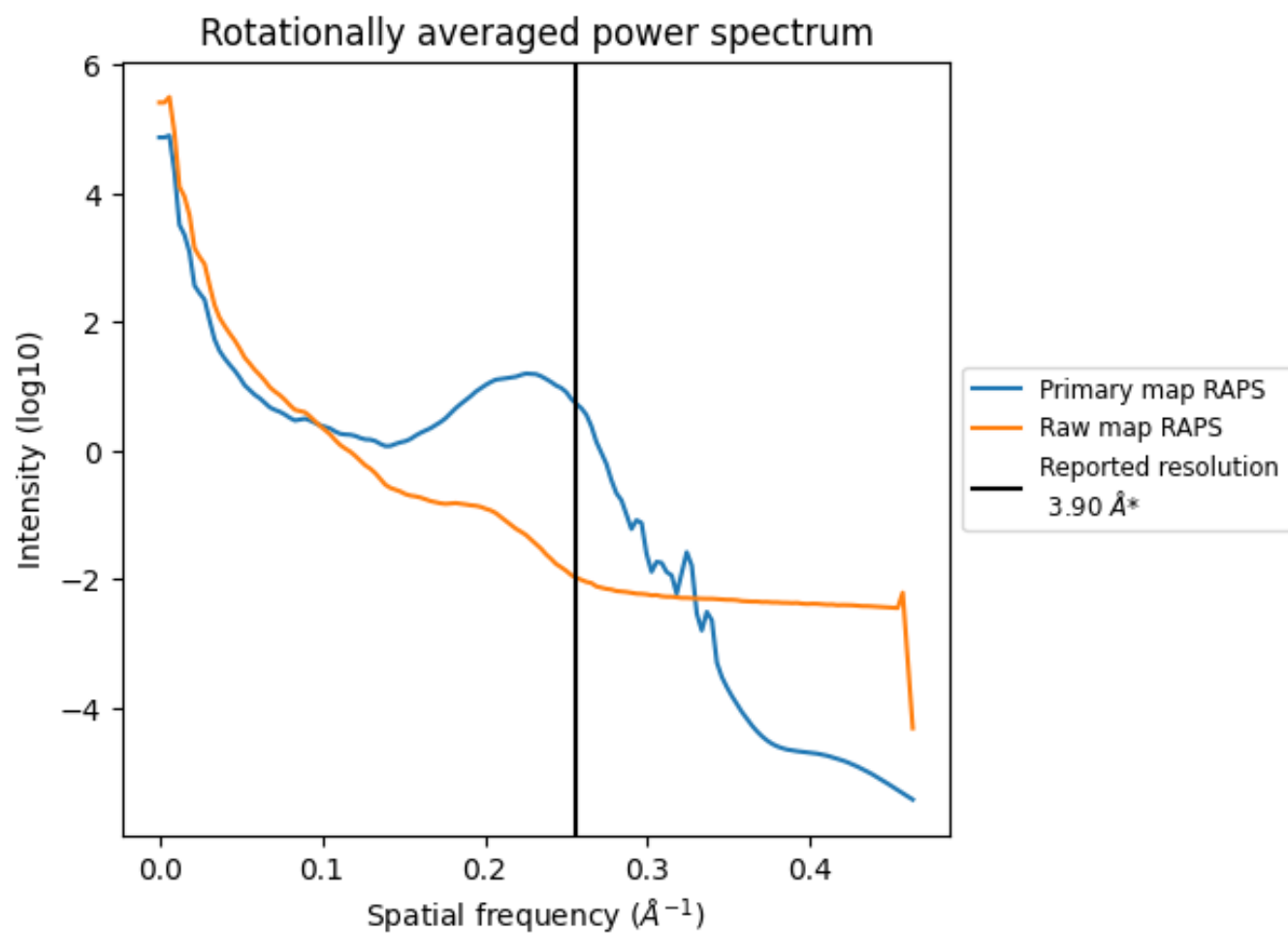
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 85 nm³; this corresponds to an approximate mass of 77 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 ⓘ

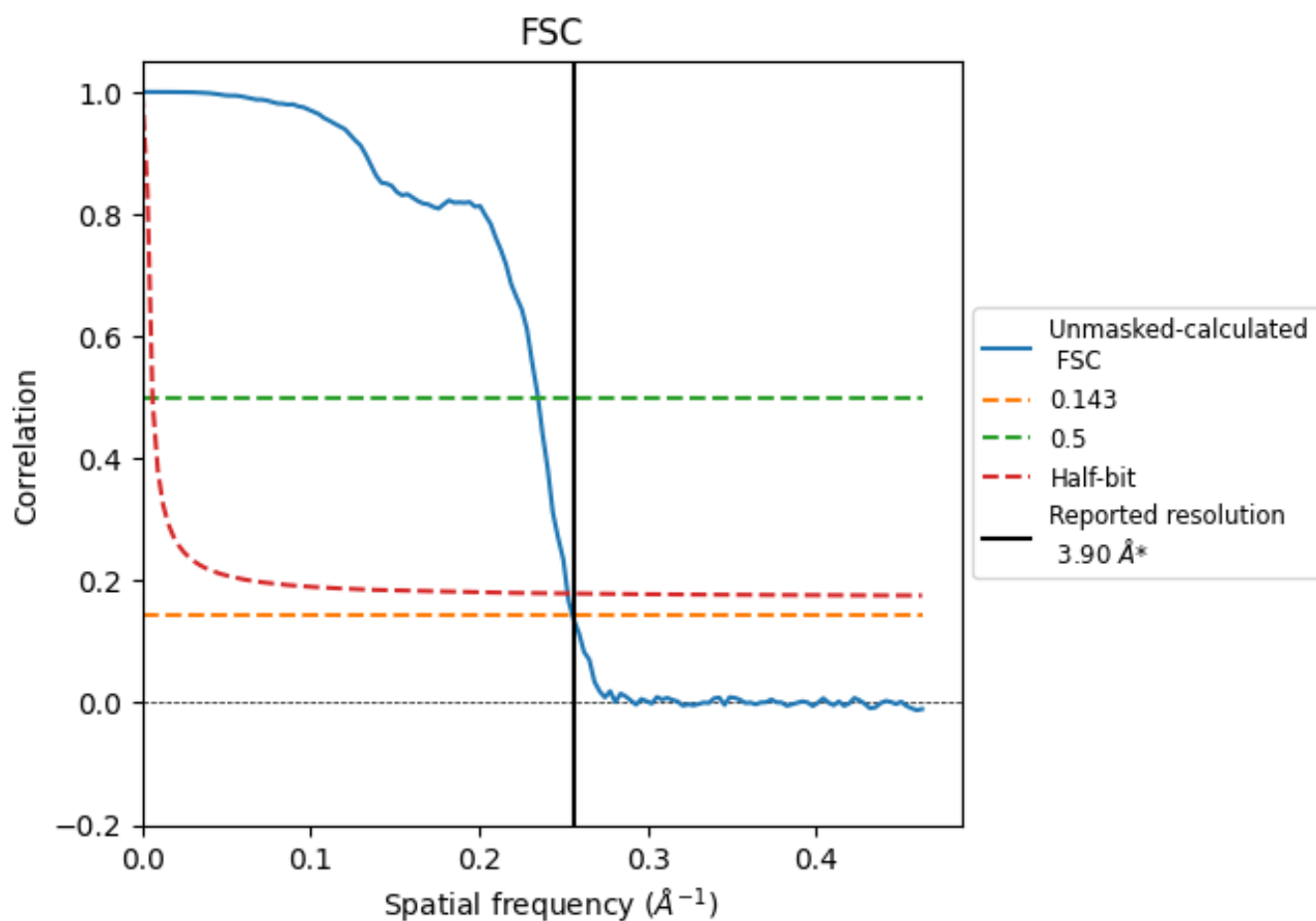


*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

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 0.256 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.92	4.26	3.96

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

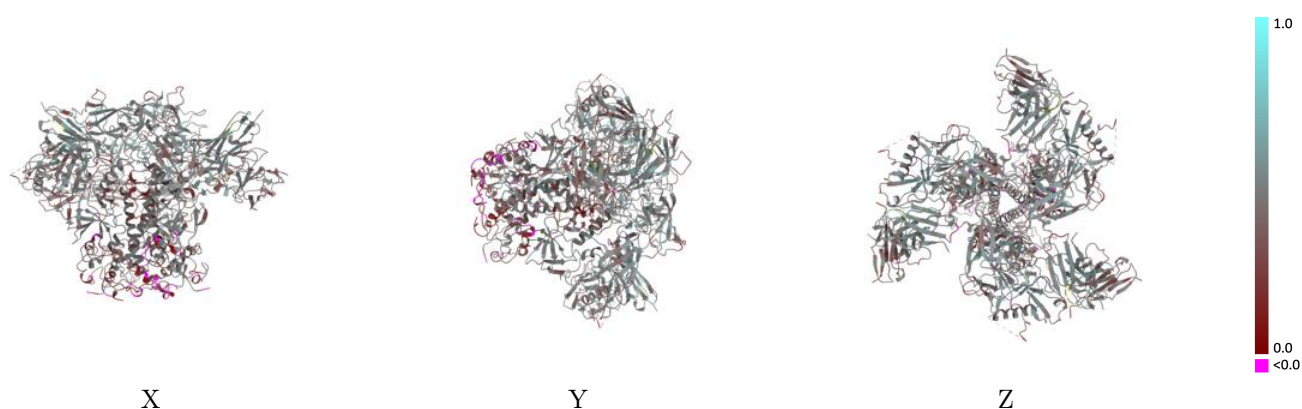
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-43228 and PDB model 8VGW. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)

This section was not generated.

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

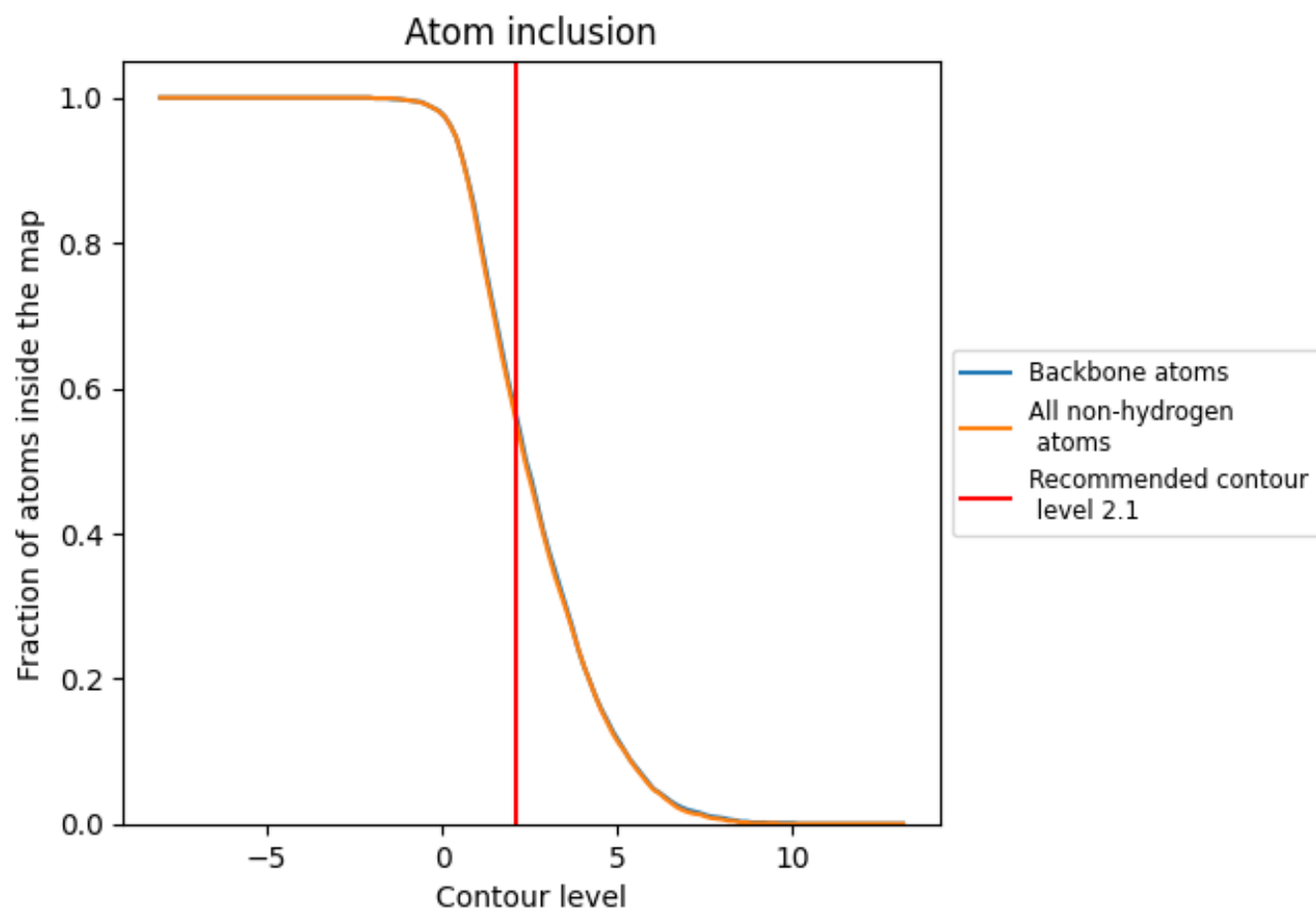


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 56% of all backbone atoms, 56% 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 (2.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5570	0.4260
A	0.6010	0.4360
B	0.4060	0.3550
C	0.6900	0.4780
D	0.4690	0.4170
E	0.6030	0.4350
F	0.4010	0.3550
G	0.6720	0.4760
H	0.4690	0.4230
I	0.6050	0.4380
J	0.4070	0.3490
K	0.6800	0.4740
L	0.4620	0.4210

1.0

0.0

<0.0