



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 03:32 PM UTC

PDB ID : 5GQH / pdb_00005gqh
EMDB ID : EMD-9535
Title : Cryo-EM structure of PaeCas3-AcrF3 complex
Authors : Zhang, X.; Ma, J.; Wang, Y.; Wang, J.
Deposited on : 2016-08-07
Resolution : 4.20 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

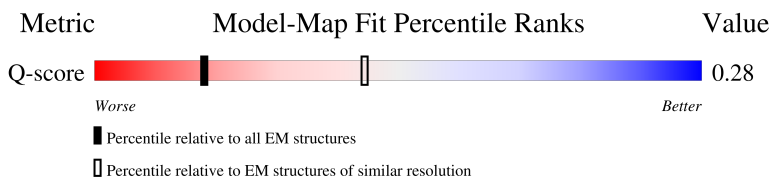
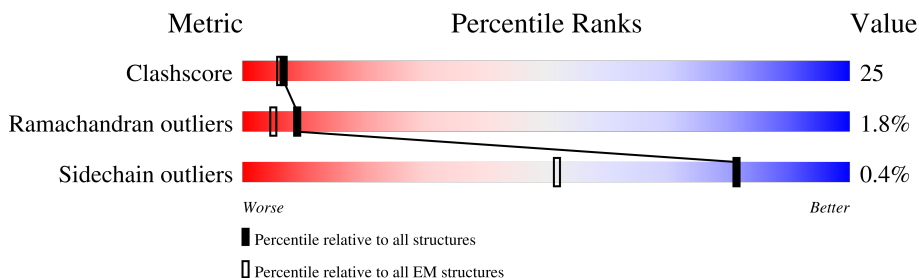
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.20 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	2937 (4.00 - 5.00)

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	1090	36% 59% 21% • 17%
2	B	139	53% 88% 9% •
2	C	139	55% 88% 9% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6641 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 CRISPR-associated nuclease/helicase Cas3 subtype I-F/YPEST.

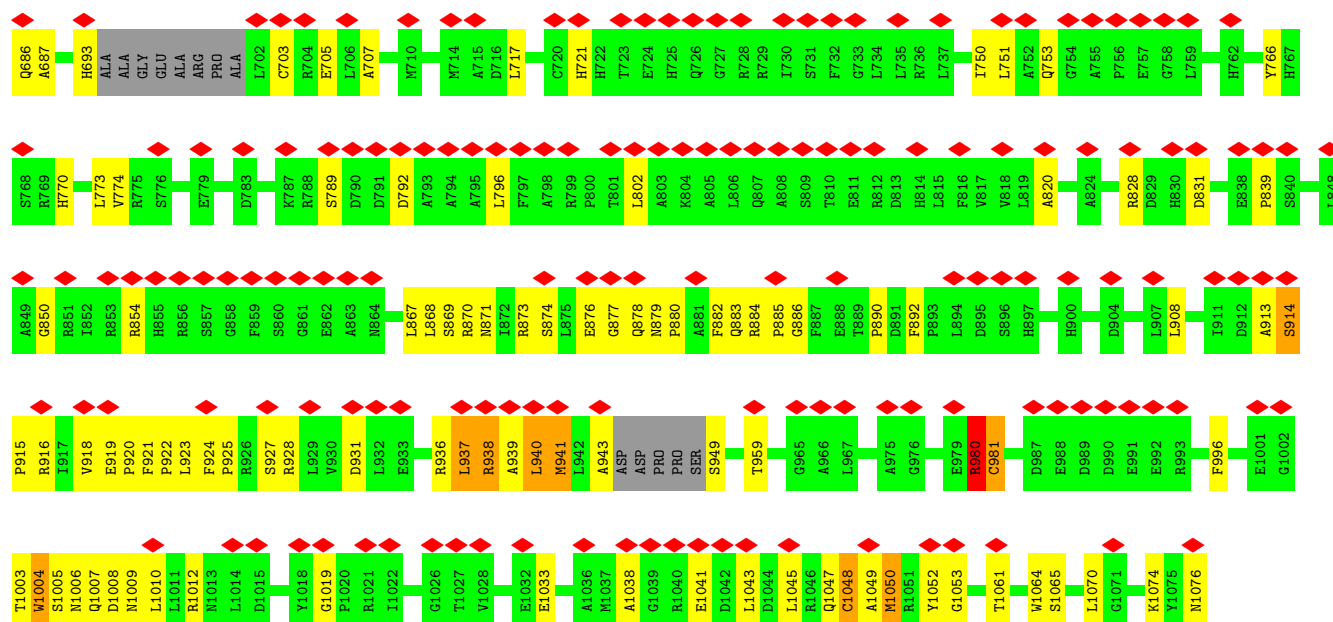
Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	901	Total	C	N	O	S	0	0
			4473	2666	904	902	1		

There are 14 discrepancies between the modelled and reference sequences:

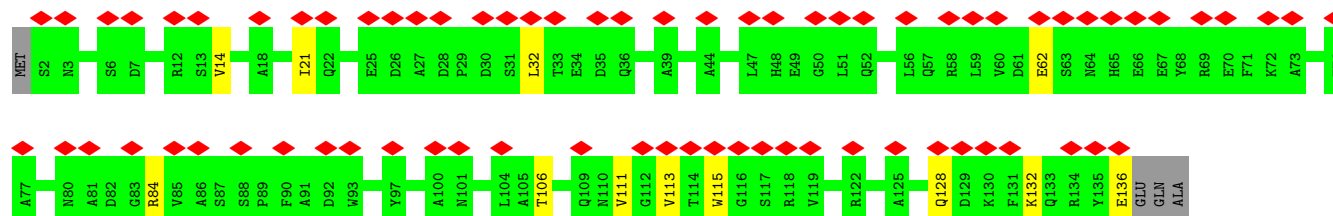
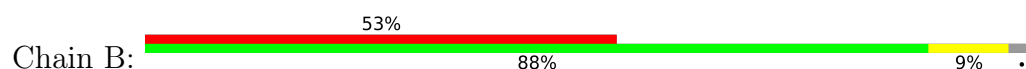
Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP Q02ML8
A	-12	GLY	-	expression tag	UNP Q02ML8
A	-11	SER	-	expression tag	UNP Q02ML8
A	-10	SER	-	expression tag	UNP Q02ML8
A	-9	HIS	-	expression tag	UNP Q02ML8
A	-8	HIS	-	expression tag	UNP Q02ML8
A	-7	HIS	-	expression tag	UNP Q02ML8
A	-6	HIS	-	expression tag	UNP Q02ML8
A	-5	HIS	-	expression tag	UNP Q02ML8
A	-4	HIS	-	expression tag	UNP Q02ML8
A	-3	SER	-	expression tag	UNP Q02ML8
A	-2	GLN	-	expression tag	UNP Q02ML8
A	-1	ASP	-	expression tag	UNP Q02ML8
A	0	PRO	-	expression tag	UNP Q02ML8

- Molecule 2 is a protein called anti-CRISPR protein 3.

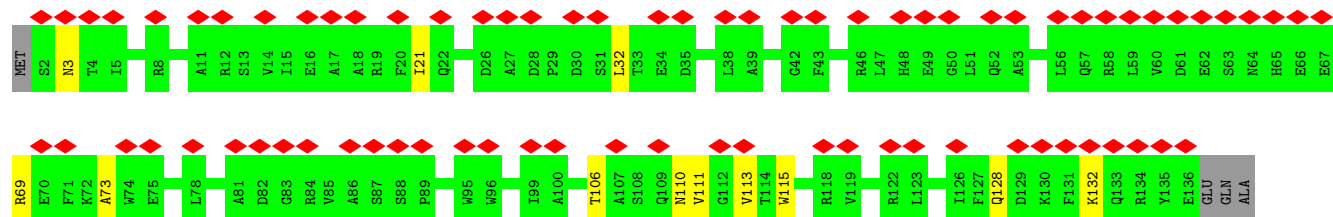
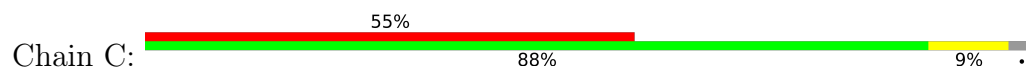
Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	135	Total	C	N	O	S	0	0
			1084	682	195	205	2		
2	C	135	Total	C	N	O	S	0	0
			1084	682	195	205	2		



• Molecule 2: anti-CRISPR protein 3



• Molecule 2: anti-CRISPR protein 3



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	78904	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	8	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.195	Depositor
Minimum map value	-0.132	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.020	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	114.799995, 114.799995, 114.799995	wwPDB
Map dimensions	140, 140, 140	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.82, 0.82, 0.82	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	1.06	17/4466 (0.4%)	1.39	73/6202 (1.2%)
2	B	0.46	2/1109 (0.2%)	0.52	0/1505
2	C	0.46	2/1109 (0.2%)	0.52	0/1505
All	All	0.90	21/6684 (0.3%)	1.18	73/9212 (0.8%)

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	5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1004	TRP	N-CA	28.79	1.81	1.46
1	A	1004	TRP	CA-C	23.77	1.81	1.52
1	A	1009	ASN	N-CA	-21.35	1.21	1.46
1	A	1050	MET	C-O	-18.02	0.99	1.23
1	A	1009	ASN	C-O	-16.95	1.04	1.24
1	A	1005	SER	N-CA	13.85	1.63	1.45
1	A	1004	TRP	C-N	10.17	1.46	1.33
1	A	576	ASP	CA-C	8.59	1.65	1.53
1	A	1008	ASP	CA-C	-8.22	1.42	1.52
1	A	1005	SER	C-O	-7.22	1.14	1.23
1	A	868	LEU	CA-C	7.15	1.62	1.52
1	A	1008	ASP	N-CA	-6.88	1.38	1.46
1	A	1006	ASN	CA-CB	-6.76	1.42	1.53
1	A	576	ASP	N-CA	6.63	1.53	1.45
2	B	115	TRP	C-O	-5.86	1.16	1.23
2	C	115	TRP	C-O	-5.86	1.16	1.23
1	A	770	HIS	C-N	5.53	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	869	SER	CA-C	5.23	1.59	1.52
2	B	32	LEU	C-O	-5.05	1.17	1.23
2	C	32	LEU	C-O	-5.04	1.17	1.23
1	A	686	GLN	CA-C	5.01	1.58	1.52

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	869	SER	N-CA-C	15.34	128.08	111.36
1	A	1050	MET	N-CA-C	-13.10	93.63	110.53
1	A	1004	TRP	N-CA-C	12.68	129.91	109.24
1	A	570	THR	N-CA-C	11.71	123.60	111.07
1	A	940	LEU	N-CA-C	11.45	127.23	110.28
1	A	940	LEU	CA-C-O	11.05	133.53	121.16
1	A	1010	LEU	N-CA-C	10.29	122.58	111.36
1	A	335	HIS	N-CA-C	9.90	123.94	108.79
1	A	370	LEU	N-CA-C	9.81	121.97	111.28
1	A	576	ASP	N-CA-C	9.07	123.95	110.52
1	A	577	GLU	N-CA-C	8.67	120.81	111.36
1	A	980	ARG	N-CA-C	8.51	120.56	111.28
1	A	868	LEU	N-CA-C	8.36	123.62	110.32
1	A	624	SER	CA-C-N	8.26	129.14	119.98
1	A	624	SER	C-N-CA	8.26	129.14	119.98
1	A	1047	GLN	CA-C-N	8.04	131.71	120.29
1	A	1047	GLN	C-N-CA	8.04	131.71	120.29
1	A	571	SER	N-CA-C	-8.04	93.67	110.80
1	A	584	ASP	N-CA-C	7.95	119.72	111.14
1	A	1005	SER	CA-C-O	-7.80	113.06	121.56
1	A	369	ARG	N-CA-C	7.59	119.55	111.28
1	A	1004	TRP	N-CA-CB	-7.37	98.42	110.43
1	A	1008	ASP	O-C-N	7.35	129.92	122.12
1	A	187	GLN	N-CA-C	7.35	119.30	111.28
1	A	118	MET	N-CA-C	7.06	118.97	111.28
1	A	333	PRO	N-CA-C	7.00	122.77	113.40
1	A	138	ASN	N-CA-C	6.91	118.82	111.28
1	A	839	PRO	O-C-N	6.72	129.45	122.93
1	A	332	PHE	CA-C-N	6.69	126.14	118.85
1	A	332	PHE	C-N-CA	6.69	126.14	118.85
1	A	128	ALA	N-CA-C	6.63	118.50	111.28
1	A	161	VAL	N-CA-C	6.63	117.41	110.72
1	A	770	HIS	CA-C-N	-6.60	112.99	119.78
1	A	770	HIS	C-N-CA	-6.60	112.99	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1050	MET	O-C-N	-6.45	114.72	122.65
1	A	575	LEU	O-C-N	-6.44	115.75	123.03
1	A	1048	CYS	N-CA-C	6.40	118.34	111.36
1	A	572	ASP	CA-C-N	6.26	130.75	122.30
1	A	572	ASP	C-N-CA	6.26	130.75	122.30
1	A	1005	SER	N-CA-C	6.25	119.32	110.50
1	A	207	PRO	N-CA-C	-6.21	104.51	110.47
1	A	1048	CYS	CA-C-N	6.16	128.54	120.28
1	A	1048	CYS	C-N-CA	6.16	128.54	120.28
1	A	1008	ASP	CA-C-O	-6.14	114.04	120.55
1	A	294	GLY	O-C-N	6.08	123.26	121.07
1	A	117	VAL	O-C-N	-5.94	115.87	121.87
1	A	680	GLN	N-CA-C	-5.91	101.29	110.10
1	A	219	SER	N-CA-C	5.81	117.61	111.28
1	A	186	SER	N-CA-C	5.76	117.56	111.28
1	A	373	GLN	N-CA-C	5.74	117.62	111.36
1	A	666	ALA	N-CA-C	-5.72	105.05	111.28
1	A	544	CYS	N-CA-C	-5.61	98.85	110.80
1	A	686	GLN	N-CA-C	-5.60	100.76	109.72
1	A	1050	MET	CA-C-O	-5.55	115.55	121.94
1	A	1047	GLN	O-C-N	5.50	128.03	122.09
1	A	182	ASP	N-CA-C	5.49	117.27	111.28
1	A	297	LEU	N-CA-C	5.46	117.31	111.36
1	A	290	LEU	N-CA-C	5.43	117.20	111.28
1	A	876	GLU	N-CA-C	-5.36	105.44	111.28
1	A	426	LYS	N-CA-C	5.35	119.81	113.28
1	A	119	ALA	N-CA-C	5.34	117.11	111.28
1	A	839	PRO	CA-C-N	5.28	131.37	123.24
1	A	839	PRO	C-N-CA	5.28	131.37	123.24
1	A	938	ARG	N-CA-C	5.23	117.06	111.36
1	A	538	GLN	N-CA-C	5.17	116.91	111.28
1	A	560	HIS	N-CA-C	-5.12	105.70	111.28
1	A	937	LEU	CA-C-N	5.10	127.53	120.29
1	A	937	LEU	C-N-CA	5.10	127.53	120.29
1	A	532	ALA	N-CA-C	5.08	116.81	111.28
1	A	1004	TRP	CA-C-N	5.07	128.89	120.94
1	A	1004	TRP	C-N-CA	5.07	128.89	120.94
1	A	1070	LEU	N-CA-C	5.05	119.45	113.28
1	A	557	ARG	N-CA-C	5.02	116.47	108.79

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	408	GLN	Peptide
1	A	575	LEU	Mainchain,Peptide
1	A	867	LEU	Mainchain
1	A	939	ALA	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	4473	0	2152	259	0
2	B	1084	0	1030	15	0
2	C	1084	0	1030	8	0
All	All	6641	0	4212	267	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 25.

All (267) 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:1004:TRP:C	1:A:1004:TRP:CA	1.82	1.53
1:A:126:GLY:HA3	1:A:153:SER:CB	1.39	1.50
1:A:1004:TRP:CA	1:A:1004:TRP:N	1.81	1.43
1:A:996:PHE:O	1:A:1007:GLN:CB	1.66	1.42
1:A:126:GLY:CA	1:A:153:SER:CA	2.01	1.36
1:A:583:ILE:CB	1:A:937:LEU:HA	1.57	1.32
1:A:126:GLY:HA3	1:A:153:SER:CA	1.60	1.29
1:A:126:GLY:CA	1:A:153:SER:HA	1.57	1.27
1:A:580:ASP:CB	1:A:586:LEU:CB	2.13	1.26
1:A:130:GLN:CB	1:A:262:PHE:CB	2.18	1.22
1:A:261:GLN:O	1:A:263:PRO:N	1.72	1.22
1:A:1012:ARG:O	1:A:1065:SER:HA	1.40	1.21
1:A:126:GLY:CA	1:A:153:SER:CB	2.19	1.20
1:A:377:LEU:CB	1:A:540:PRO:CB	2.20	1.18
1:A:594:HIS:CB	1:A:621:ALA:HB1	1.73	1.18
1:A:890:PRO:CB	2:B:62:GLU:HB3	1.76	1.15
1:A:146:ALA:O	1:A:147:TYR:O	1.61	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:ASP:CB	1:A:854:ARG:CB	2.24	1.14
1:A:402:ALA:HB1	1:A:436:ALA:CB	1.77	1.14
1:A:180:THR:CB	1:A:181:GLY:HA2	1.75	1.13
1:A:126:GLY:HA2	1:A:153:SER:HA	1.13	1.12
1:A:583:ILE:CB	1:A:937:LEU:CA	2.30	1.08
1:A:580:ASP:CB	1:A:586:LEU:CA	2.32	1.07
1:A:129:SER:CB	1:A:259:CYS:O	2.03	1.04
1:A:126:GLY:HA2	1:A:153:SER:CA	1.77	1.04
1:A:394:TRP:CB	1:A:649:GLU:HA	1.88	1.03
1:A:338:THR:CB	1:A:344:LYS:O	2.07	1.03
1:A:402:ALA:HB1	1:A:436:ALA:HB1	1.01	0.99
1:A:643:ARG:HA	1:A:655:SER:O	1.63	0.99
1:A:120:ALA:CB	1:A:356:VAL:CB	2.41	0.98
1:A:402:ALA:CB	1:A:436:ALA:HB1	1.93	0.98
1:A:580:ASP:CB	1:A:586:LEU:HA	1.95	0.95
1:A:405:CYS:O	1:A:409:ALA:HB2	1.66	0.94
1:A:597:GLY:O	1:A:600:GLY:N	2.00	0.94
1:A:338:THR:CB	1:A:345:LEU:HA	1.98	0.94
1:A:831:ASP:CA	1:A:854:ARG:CB	2.47	0.93
1:A:159:ALA:HB2	1:A:188:LEU:CB	2.00	0.90
1:A:996:PHE:C	1:A:1007:GLN:CB	2.44	0.90
1:A:883:GLN:HA	1:A:886:GLY:HA3	1.54	0.90
1:A:131:ALA:H	1:A:259:CYS:HA	1.37	0.90
1:A:377:LEU:N	1:A:540:PRO:CB	2.34	0.90
1:A:423:GLY:HA3	1:A:850:GLY:HA2	1.53	0.88
1:A:297:LEU:O	1:A:299:HIS:N	2.06	0.87
1:A:417:LEU:HA	1:A:605:LEU:O	1.74	0.87
1:A:375:PRO:CB	1:A:538:GLN:O	2.23	0.86
1:A:159:ALA:CB	1:A:188:LEU:CB	2.54	0.86
1:A:120:ALA:HB2	1:A:356:VAL:CB	2.06	0.85
1:A:980:ARG:O	1:A:981:CYS:CB	2.23	0.85
1:A:405:CYS:O	1:A:409:ALA:CB	2.24	0.85
1:A:873:ARG:O	1:A:878:GLN:N	2.11	0.83
1:A:583:ILE:CB	1:A:936:ARG:O	2.27	0.83
1:A:120:ALA:HB1	1:A:356:VAL:CB	2.08	0.83
1:A:402:ALA:CB	1:A:436:ALA:CB	2.55	0.81
1:A:883:GLN:HA	1:A:886:GLY:CA	2.10	0.81
1:A:583:ILE:C	1:A:940:LEU:CB	2.54	0.81
1:A:418:ASN:O	1:A:606:SER:HA	1.80	0.81
1:A:913:ALA:O	1:A:915:PRO:N	2.13	0.81
1:A:918:VAL:HA	1:A:919:GLU:CB	2.09	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:GLY:CA	1:A:153:SER:N	2.42	0.81
1:A:471:ASP:O	1:A:474:LEU:O	1.98	0.80
1:A:773:LEU:HD12	1:A:774:VAL:N	1.95	0.80
1:A:477:LEU:N	1:A:543:ALA:O	2.14	0.80
1:A:180:THR:CB	1:A:181:GLY:CA	2.56	0.79
1:A:1043:LEU:CB	1:A:1048:CYS:CB	2.61	0.79
1:A:873:ARG:O	1:A:877:GLY:N	2.15	0.79
1:A:212:ALA:HB2	1:A:298:LEU:CB	2.13	0.79
1:A:582:ASP:O	1:A:583:ILE:C	2.24	0.78
1:A:924:PHE:CB	1:A:927:SER:CB	2.64	0.76
1:A:409:ALA:O	1:A:413:GLY:N	2.17	0.76
1:A:1004:TRP:C	1:A:1004:TRP:HA	2.09	0.76
1:A:1076:ASN:O	2:C:73:ALA:HB2	1.85	0.76
1:A:261:GLN:C	1:A:263:PRO:N	2.41	0.76
1:A:1004:TRP:N	1:A:1004:TRP:CB	2.49	0.76
1:A:363:LEU:O	1:A:367:LEU:HD13	1.85	0.76
1:A:580:ASP:O	1:A:585:ASP:CB	2.33	0.75
1:A:949:SER:HA	2:B:84:ARG:NH1	2.01	0.75
1:A:871:ASN:O	1:A:874:SER:N	2.18	0.75
1:A:423:GLY:CA	1:A:850:GLY:CA	2.64	0.75
1:A:214:GLY:O	1:A:218:VAL:N	2.21	0.74
1:A:236:ARG:O	1:A:237:LEU:O	2.06	0.74
1:A:831:ASP:HA	1:A:854:ARG:CB	2.17	0.73
1:A:882:PHE:C	1:A:886:GLY:HA3	2.13	0.73
1:A:126:GLY:HA3	1:A:153:SER:N	2.00	0.73
1:A:924:PHE:O	1:A:927:SER:N	2.20	0.73
1:A:414:PHE:O	1:A:602:ARG:CB	2.37	0.73
1:A:1038:ALA:O	1:A:1041:GLU:O	2.07	0.73
1:A:582:ASP:O	1:A:584:ASP:N	2.22	0.73
1:A:146:ALA:O	1:A:147:TYR:C	2.31	0.73
1:A:913:ALA:O	1:A:914:SER:C	2.31	0.72
1:A:687:ALA:HB3	1:A:908:LEU:O	1.89	0.72
1:A:423:GLY:HA3	1:A:850:GLY:CA	2.19	0.72
1:A:583:ILE:CB	1:A:937:LEU:C	2.63	0.72
1:A:594:HIS:CA	1:A:621:ALA:HB1	2.19	0.72
1:A:193:ARG:N	1:A:194:GLN:HA	2.05	0.72
1:A:131:ALA:HB3	1:A:259:CYS:CB	2.19	0.71
1:A:693:HIS:O	1:A:705:GLU:CB	2.38	0.71
1:A:297:LEU:O	1:A:298:LEU:C	2.32	0.71
1:A:703:CYS:O	1:A:707:ALA:N	2.23	0.71
1:A:377:LEU:CA	1:A:540:PRO:CB	2.67	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:883:GLN:CA	1:A:886:GLY:HA3	2.20	0.71
1:A:583:ILE:CA	1:A:937:LEU:HA	2.21	0.71
1:A:980:ARG:O	1:A:981:CYS:HB2	1.90	0.71
1:A:261:GLN:O	1:A:262:PHE:C	2.33	0.70
1:A:236:ARG:O	1:A:237:LEU:C	2.35	0.70
1:A:583:ILE:CB	1:A:940:LEU:CB	2.69	0.70
1:A:1076:ASN:C	2:C:69:ARG:HG3	2.17	0.70
1:A:594:HIS:CB	1:A:621:ALA:CB	2.64	0.70
1:A:883:GLN:HA	1:A:886:GLY:N	2.07	0.69
1:A:418:ASN:O	1:A:606:SER:CA	2.39	0.69
1:A:126:GLY:HA2	1:A:153:SER:N	2.05	0.68
1:A:643:ARG:CA	1:A:655:SER:O	2.41	0.68
1:A:174:LEU:O	1:A:278:ALA:HA	1.94	0.68
1:A:126:GLY:N	1:A:153:SER:CB	2.56	0.67
1:A:684:ARG:O	1:A:685:ARG:CB	2.41	0.67
1:A:773:LEU:HD12	1:A:773:LEU:C	2.19	0.67
1:A:263:PRO:CB	1:A:264:HIS:HA	2.24	0.67
1:A:130:GLN:N	1:A:259:CYS:O	2.28	0.67
1:A:597:GLY:O	1:A:600:GLY:CA	2.43	0.66
1:A:879:ASN:HA	1:A:880:PRO:C	2.21	0.66
1:A:377:LEU:N	1:A:444:ALA:O	2.24	0.66
1:A:890:PRO:N	2:B:62:GLU:OE1	2.28	0.66
1:A:261:GLN:O	1:A:263:PRO:CA	2.43	0.66
1:A:261:GLN:O	1:A:263:PRO:CB	2.45	0.65
1:A:228:HIS:CB	2:B:136:GLU:H	2.09	0.65
1:A:126:GLY:N	1:A:153:SER:HA	2.09	0.65
1:A:949:SER:HA	2:B:84:ARG:HH11	1.60	0.65
1:A:131:ALA:N	1:A:259:CYS:HA	2.10	0.65
1:A:1004:TRP:C	1:A:1004:TRP:CB	2.68	0.65
1:A:131:ALA:CB	1:A:259:CYS:HA	2.27	0.65
1:A:423:GLY:CA	1:A:850:GLY:HA3	2.27	0.65
1:A:890:PRO:CB	2:B:62:GLU:CB	2.66	0.65
1:A:171:LEU:O	1:A:281:ALA:HB1	1.97	0.64
1:A:1003:THR:C	1:A:1004:TRP:CA	2.68	0.64
1:A:938:ARG:HA	1:A:940:LEU:O	1.96	0.63
2:B:14:VAL:HG22	2:C:110:ASN:OD1	1.98	0.63
1:A:940:LEU:O	1:A:941:MET:CB	2.45	0.63
1:A:130:GLN:O	1:A:131:ALA:C	2.40	0.63
1:A:175:ALA:HA	1:A:278:ALA:HA	1.81	0.62
1:A:980:ARG:O	1:A:981:CYS:HB3	1.98	0.62
1:A:938:ARG:O	1:A:943:ALA:HB2	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:PHE:O	1:A:644:CYS:HA	2.00	0.61
1:A:175:ALA:O	1:A:278:ALA:HB1	1.99	0.61
1:A:559:GLY:O	1:A:562:ILE:N	2.34	0.61
1:A:129:SER:CA	1:A:259:CYS:O	2.48	0.61
1:A:572:ASP:HA	1:A:601:SER:CB	2.31	0.61
1:A:130:GLN:O	1:A:133:GLN:N	2.34	0.61
1:A:870:ARG:CB	1:A:874:SER:CB	2.79	0.60
1:A:882:PHE:O	1:A:886:GLY:HA3	2.00	0.60
1:A:648:ASP:CB	1:A:651:SER:O	2.50	0.60
1:A:129:SER:CB	1:A:259:CYS:C	2.75	0.59
1:A:175:ALA:HB2	1:A:281:ALA:HB3	1.84	0.59
1:A:129:SER:C	1:A:259:CYS:O	2.45	0.59
1:A:789:SER:O	1:A:792:ASP:CB	2.51	0.58
1:A:228:HIS:CB	2:B:136:GLU:N	2.66	0.58
1:A:583:ILE:O	1:A:940:LEU:CB	2.52	0.58
1:A:399:TYR:O	1:A:402:ALA:HB3	2.03	0.58
2:B:128:GLN:O	2:B:132:LYS:HG3	2.03	0.58
2:C:128:GLN:O	2:C:132:LYS:HG3	2.03	0.58
1:A:130:GLN:CB	1:A:262:PHE:CA	2.82	0.58
1:A:416:GLY:O	1:A:605:LEU:N	2.36	0.57
1:A:131:ALA:CB	1:A:259:CYS:CB	2.83	0.57
1:A:773:LEU:HB3	1:A:1052:TYR:O	2.03	0.57
1:A:583:ILE:CB	1:A:937:LEU:O	2.53	0.57
1:A:796:LEU:O	1:A:802:LEU:CB	2.53	0.57
1:A:874:SER:C	1:A:877:GLY:H	2.13	0.56
1:A:913:ALA:O	1:A:916:ARG:N	2.38	0.56
1:A:1012:ARG:O	1:A:1065:SER:CA	2.33	0.56
1:A:233:SER:CB	2:B:136:GLU:OE1	2.54	0.56
1:A:263:PRO:CB	1:A:264:HIS:CA	2.85	0.55
1:A:607:SER:CB	1:A:610:LEU:CB	2.86	0.54
1:A:883:GLN:N	1:A:886:GLY:HA3	2.23	0.54
1:A:923:LEU:CB	1:A:931:ASP:CB	2.85	0.54
1:A:297:LEU:C	1:A:299:HIS:N	2.66	0.54
1:A:446:PHE:O	1:A:541:ILE:HA	2.08	0.54
1:A:583:ILE:CB	1:A:936:ARG:C	2.81	0.53
1:A:773:LEU:CD1	1:A:1033:GLU:CB	2.87	0.53
1:A:145:ASP:CB	1:A:146:ALA:CA	2.85	0.53
1:A:402:ALA:CB	1:A:436:ALA:HB2	2.37	0.53
1:A:646:TRP:O	1:A:652:SER:HA	2.08	0.53
1:A:131:ALA:CB	1:A:259:CYS:CA	2.87	0.53
1:A:477:LEU:O	1:A:544:CYS:HA	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ALA:HB1	1:A:188:LEU:CB	2.39	0.53
1:A:370:LEU:O	1:A:374:LEU:N	2.34	0.52
1:A:159:ALA:HB1	1:A:188:LEU:CA	2.39	0.52
1:A:423:GLY:HA2	1:A:850:GLY:CA	2.39	0.52
1:A:918:VAL:CA	1:A:919:GLU:CB	2.85	0.52
1:A:131:ALA:HB2	1:A:259:CYS:CA	2.39	0.52
1:A:131:ALA:HB2	1:A:259:CYS:HA	1.92	0.52
1:A:175:ALA:HB2	1:A:281:ALA:CB	2.40	0.52
1:A:664:SER:O	1:A:668:ALA:N	2.44	0.51
1:A:883:GLN:CA	1:A:886:GLY:N	2.73	0.51
1:A:126:GLY:H	1:A:153:SER:CB	2.23	0.51
1:A:924:PHE:O	1:A:925:PRO:C	2.53	0.51
1:A:159:ALA:CB	1:A:188:LEU:CA	2.88	0.51
1:A:884:ARG:N	1:A:885:PRO:CA	2.73	0.51
1:A:928:ARG:CB	1:A:931:ASP:CB	2.90	0.50
1:A:145:ASP:CB	1:A:146:ALA:HA	2.41	0.50
1:A:206:LEU:CB	1:A:210:ALA:HB3	2.41	0.50
1:A:884:ARG:N	1:A:885:PRO:HA	2.25	0.50
1:A:477:LEU:CB	1:A:544:CYS:CB	2.90	0.49
1:A:582:ASP:C	1:A:584:ASP:N	2.70	0.49
1:A:750:ILE:O	1:A:753:GLN:O	2.30	0.49
2:C:21:ILE:HD11	2:C:106:THR:HG22	1.95	0.49
1:A:145:ASP:CB	1:A:146:ALA:C	2.85	0.49
1:A:918:VAL:CB	1:A:919:GLU:C	2.85	0.49
1:A:717:LEU:O	1:A:721:HIS:N	2.45	0.49
1:A:418:ASN:C	1:A:606:SER:HA	2.39	0.48
1:A:874:SER:O	1:A:877:GLY:HA2	2.12	0.48
1:A:159:ALA:HB1	1:A:188:LEU:N	2.28	0.48
1:A:175:ALA:HA	1:A:278:ALA:CA	2.44	0.48
1:A:572:ASP:CB	1:A:602:ARG:O	2.62	0.48
1:A:475:ALA:HA	1:A:513:HIS:O	2.14	0.47
1:A:773:LEU:HD13	1:A:1033:GLU:CB	2.44	0.47
1:A:883:GLN:CA	1:A:886:GLY:CA	2.86	0.47
2:B:21:ILE:HD11	2:B:106:THR:HG22	1.95	0.47
1:A:766:TYR:O	1:A:820:ALA:HA	2.15	0.47
1:A:163:PRO:HA	1:A:164:GLY:HA2	1.75	0.46
1:A:399:TYR:O	1:A:402:ALA:CB	2.63	0.46
1:A:175:ALA:CA	1:A:281:ALA:HB3	2.45	0.46
1:A:297:LEU:O	1:A:300:ASP:N	2.40	0.46
1:A:130:GLN:CB	1:A:262:PHE:C	2.90	0.45
1:A:678:LEU:CB	1:A:914:SER:CB	2.94	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1061:THR:CB	2:C:3:ASN:ND2	2.80	0.45
1:A:147:TYR:O	1:A:148:ARG:C	2.58	0.45
1:A:423:GLY:HA2	1:A:850:GLY:HA3	1.98	0.45
1:A:773:LEU:HD12	1:A:774:VAL:CA	2.47	0.45
1:A:890:PRO:CB	2:B:62:GLU:OE1	2.64	0.45
1:A:773:LEU:CD1	1:A:774:VAL:N	2.73	0.44
1:A:871:ASN:N	1:A:874:SER:CB	2.80	0.44
1:A:913:ALA:C	1:A:915:PRO:N	2.75	0.44
1:A:597:GLY:C	1:A:600:GLY:H	2.16	0.44
1:A:443:GLY:HA2	1:A:444:ALA:HA	1.47	0.44
1:A:871:ASN:C	1:A:874:SER:H	2.26	0.44
1:A:643:ARG:HA	1:A:656:ALA:HA	1.99	0.44
1:A:1012:ARG:N	1:A:1064:TRP:O	2.44	0.44
1:A:175:ALA:CB	1:A:281:ALA:HB3	2.48	0.43
1:A:921:PHE:HA	1:A:922:PRO:HA	1.83	0.43
1:A:175:ALA:HA	1:A:278:ALA:O	2.18	0.43
1:A:597:GLY:O	1:A:600:GLY:HA2	2.17	0.43
1:A:687:ALA:CB	1:A:908:LEU:O	2.61	0.43
1:A:179:GLU:HA	1:A:180:THR:HA	1.71	0.42
1:A:378:ALA:HA	1:A:444:ALA:HB2	2.01	0.42
1:A:831:ASP:N	1:A:854:ARG:CB	2.82	0.42
1:A:1049:ALA:O	1:A:1053:GLY:N	2.52	0.42
1:A:299:HIS:O	1:A:300:ASP:C	2.62	0.42
1:A:235:ALA:HA	1:A:959:THR:CB	2.50	0.42
1:A:233:SER:CB	2:B:136:GLU:CD	2.93	0.42
2:B:111:VAL:HG23	2:B:113:VAL:HG13	2.02	0.42
1:A:678:LEU:O	1:A:680:GLN:O	2.38	0.41
1:A:521:GLY:HA3	1:A:524:ARG:CB	2.51	0.41
1:A:594:HIS:HA	1:A:621:ALA:HB1	2.00	0.41
1:A:773:LEU:HD11	1:A:1033:GLU:CB	2.51	0.41
1:A:1045:LEU:O	1:A:1049:ALA:HB3	2.20	0.41
1:A:207:PRO:O	1:A:210:ALA:HB3	2.21	0.41
1:A:883:GLN:C	1:A:886:GLY:N	2.79	0.41
1:A:236:ARG:C	1:A:237:LEU:O	2.64	0.41
1:A:555:SER:CB	1:A:561:GLN:CB	2.99	0.41
2:C:111:VAL:HG23	2:C:113:VAL:HG13	2.02	0.41
1:A:215:TRP:O	1:A:219:SER:N	2.47	0.41
1:A:703:CYS:O	1:A:707:ALA:HB2	2.21	0.41
2:B:14:VAL:HG22	2:C:110:ASN:CG	2.45	0.41
1:A:416:GLY:O	1:A:604:LEU:HA	2.21	0.40
1:A:645:ALA:HA	1:A:654:SER:HA	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:664:SER:O	1:A:667:HIS:CB	2.68	0.40
1:A:439:ASP:HA	1:A:440:PRO:HA	1.84	0.40
1:A:751:LEU:O	1:A:1019:GLY:N	2.54	0.40
1:A:1048:CYS:C	1:A:1050:MET:O	2.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	885/1090 (81%)	837 (95%)	27 (3%)	21 (2%)	4	28
2	B	133/139 (96%)	129 (97%)	4 (3%)	0	100	100
2	C	133/139 (96%)	129 (97%)	4 (3%)	0	100	100
All	All	1151/1368 (84%)	1095 (95%)	35 (3%)	21 (2%)	9	33

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	ASP
1	A	147	TYR
1	A	197	PRO
1	A	262	PHE
1	A	263	PRO
1	A	298	LEU
1	A	342	SER
1	A	544	CYS
1	A	685	ARG
1	A	828	ARG
1	A	892	PHE
1	A	981	CYS
1	A	1074	LYS

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Mol	Chain	Res	Type
1	A	339	ASP
1	A	941	MET
1	A	163	PRO
1	A	344	LYS
1	A	914	SER
1	A	920	PRO
1	A	237	LEU
1	A	583	ILE

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	7/898 (1%)	6 (86%)	1 (14%)	3	16
2	B	109/112 (97%)	109 (100%)	0	100	100
2	C	109/112 (97%)	109 (100%)	0	100	100
All	All	225/1122 (20%)	224 (100%)	1 (0%)	81	82

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	980	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

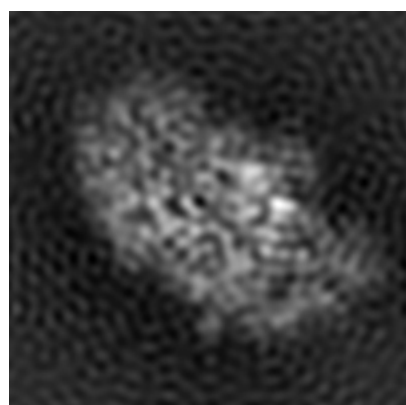
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9535. These allow visual inspection of the internal detail of the map and identification of artifacts.

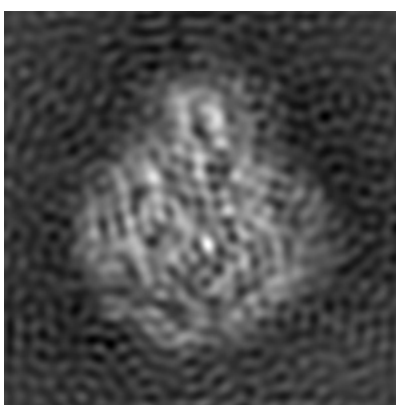
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

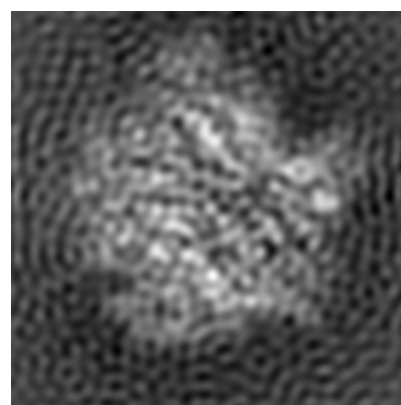
6.1.1 Primary map



X



Y

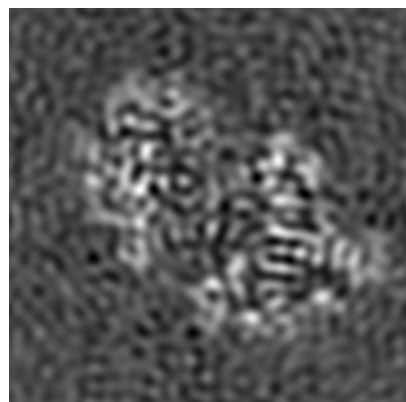


Z

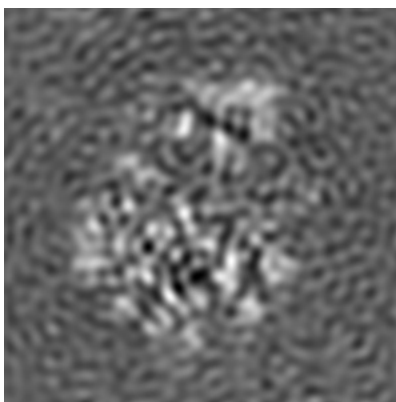
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

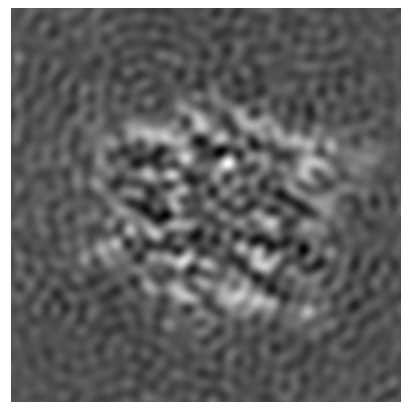
6.2.1 Primary map



X Index: 70



Y Index: 70

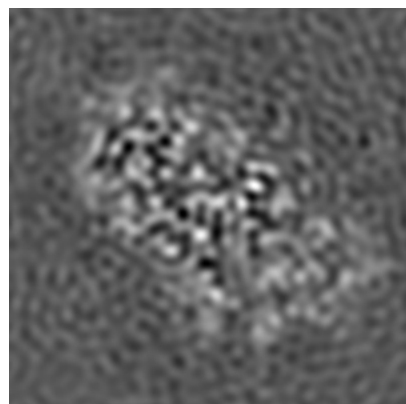


Z Index: 70

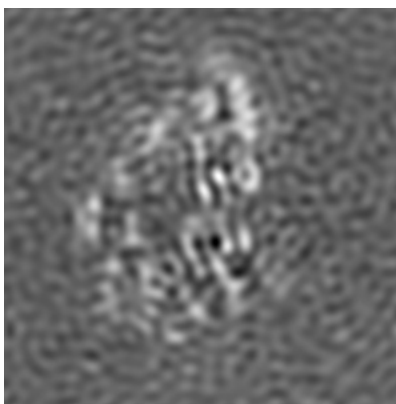
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

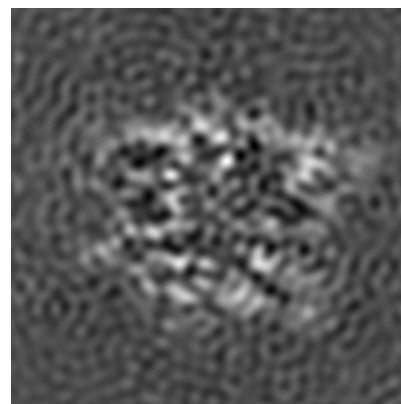
6.3.1 Primary map



X Index: 56



Y Index: 84

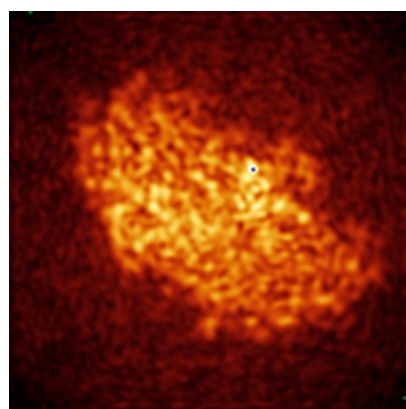


Z Index: 71

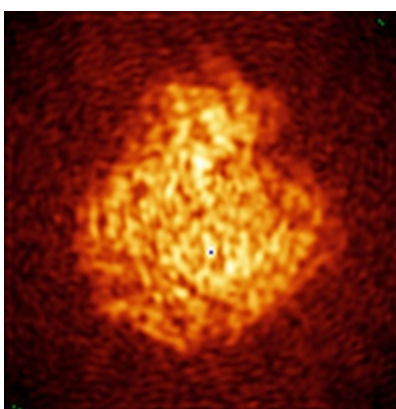
The images above show the largest variance slices of the map in three orthogonal directions.

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

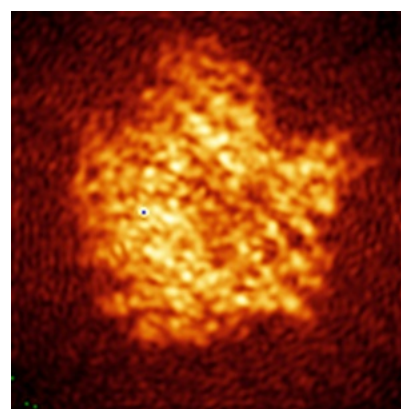
6.4.1 Primary map



X



Y

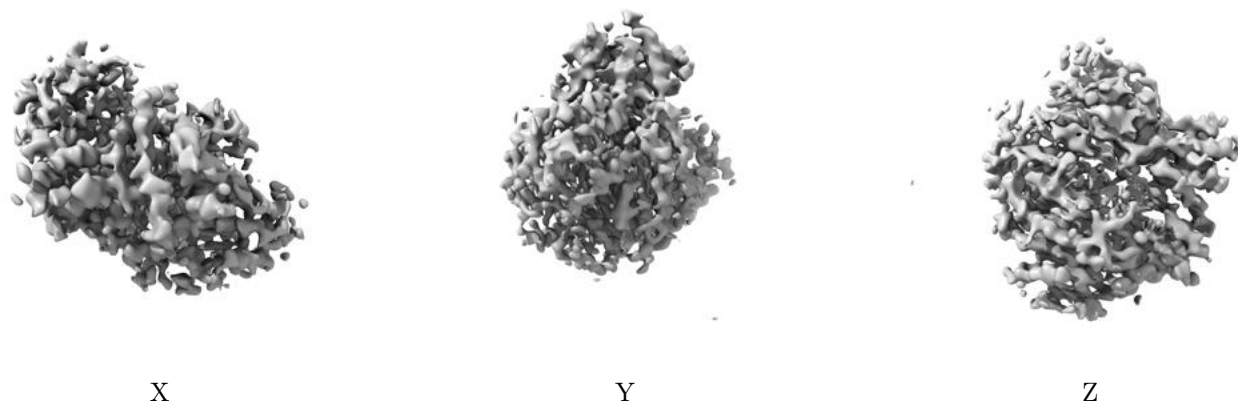


Z

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

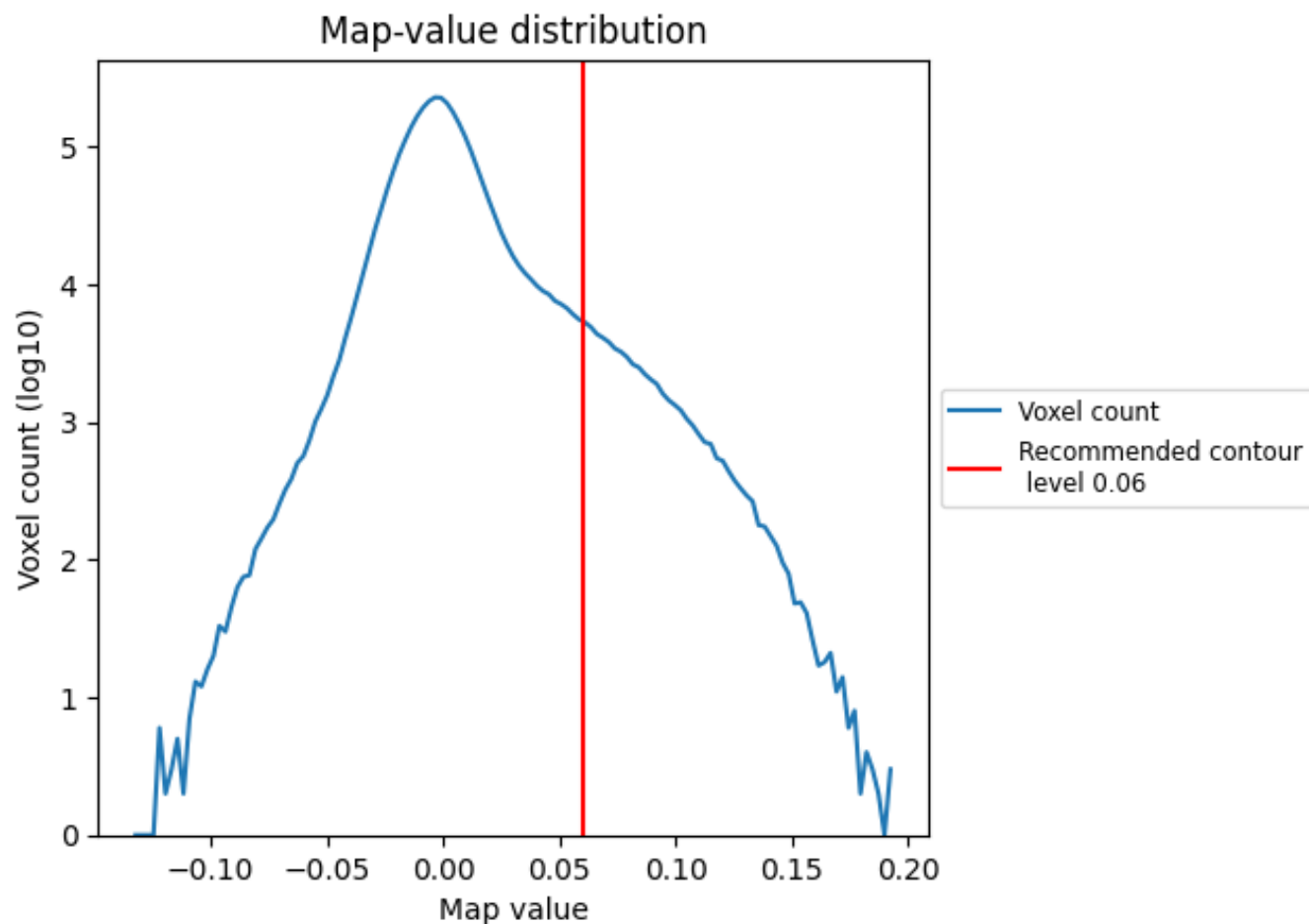
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

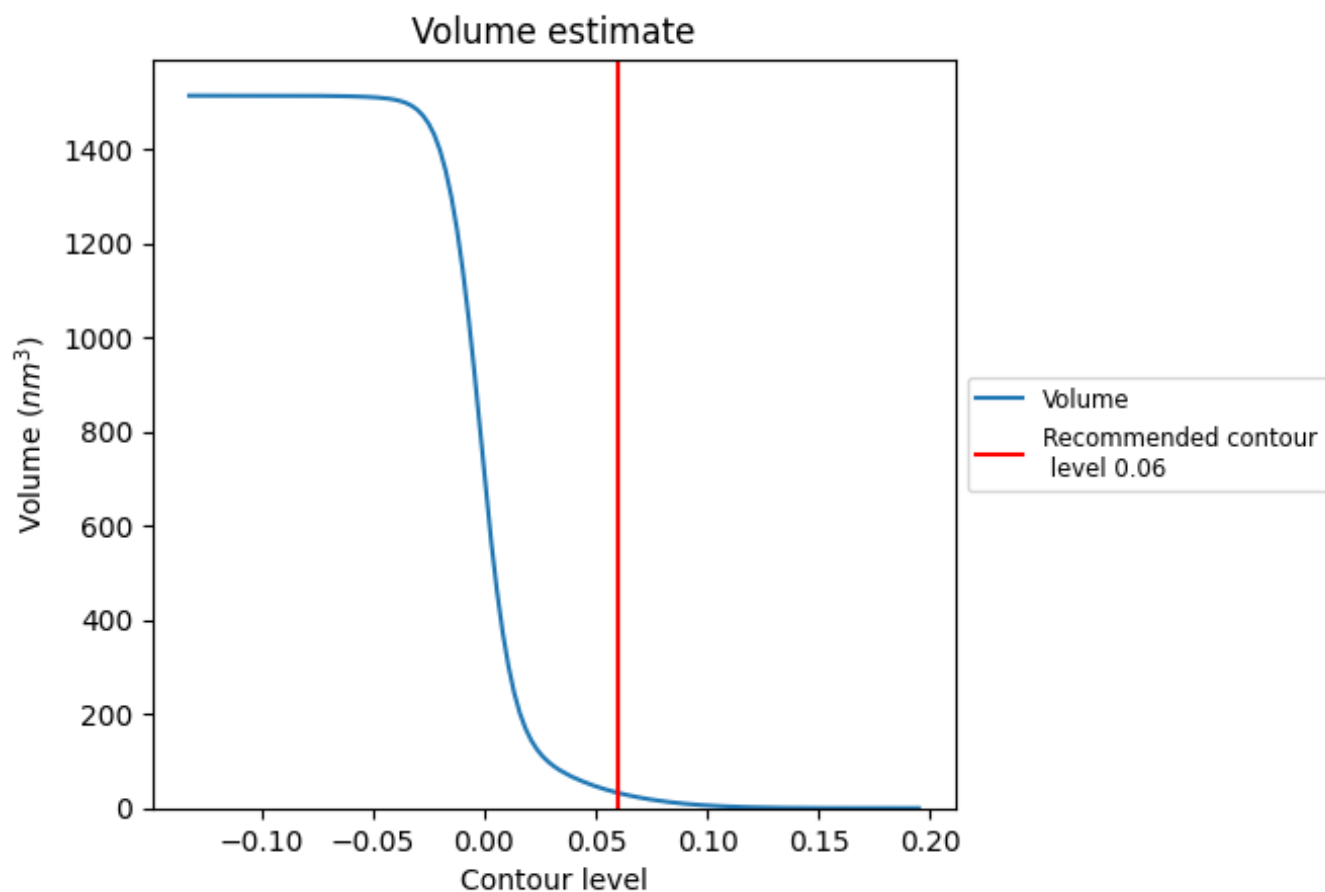
This section contains the results of statistical analysis of the map.

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.

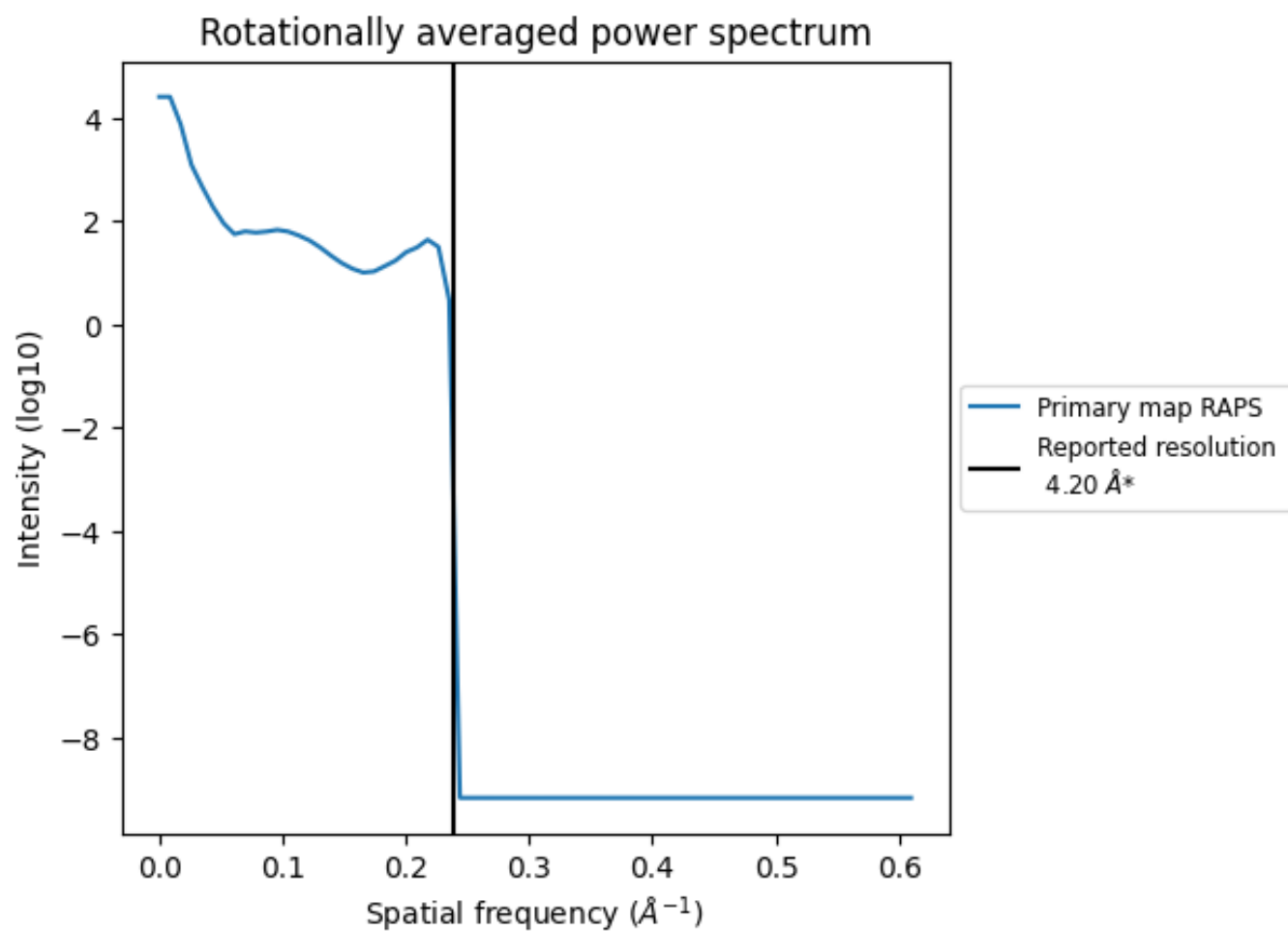
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 33 nm³; this corresponds to an approximate mass of 29 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 0.238 $\text{\AA}^{-1}$

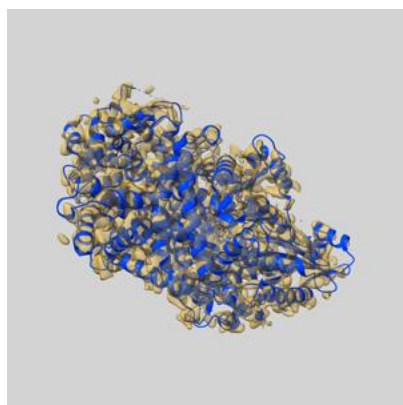
8 Fourier-Shell correlation ⓘ

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

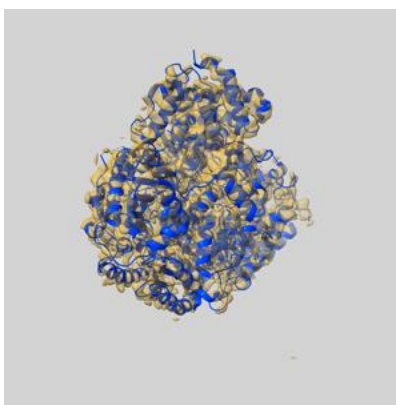
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-9535 and PDB model 5GQH. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

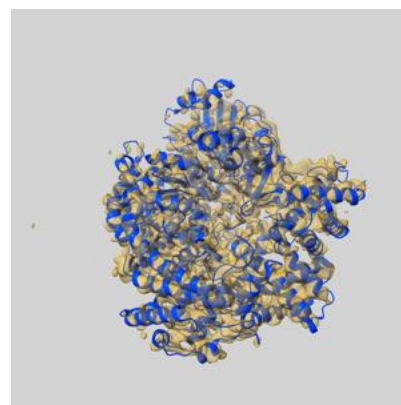
9.1 Map-model overlay [i](#)



X



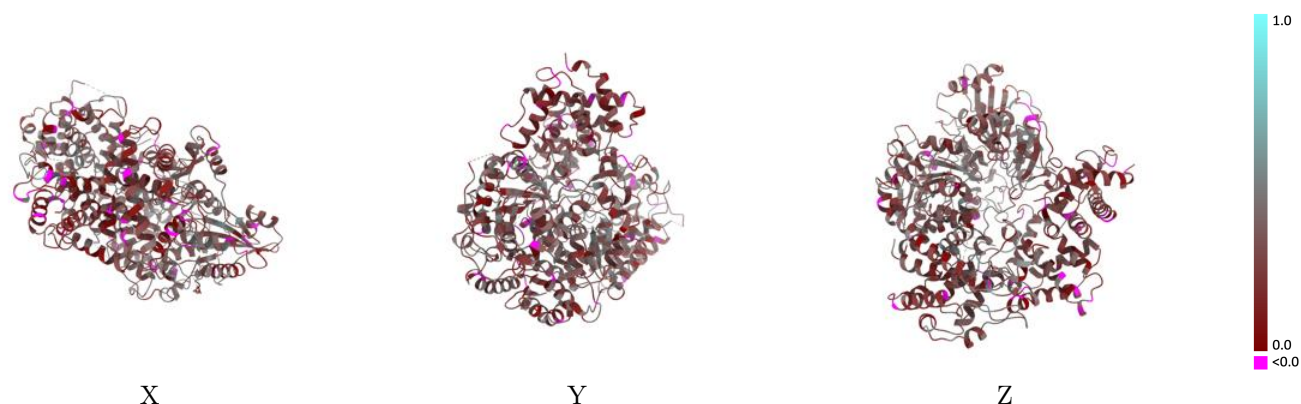
Y



Z

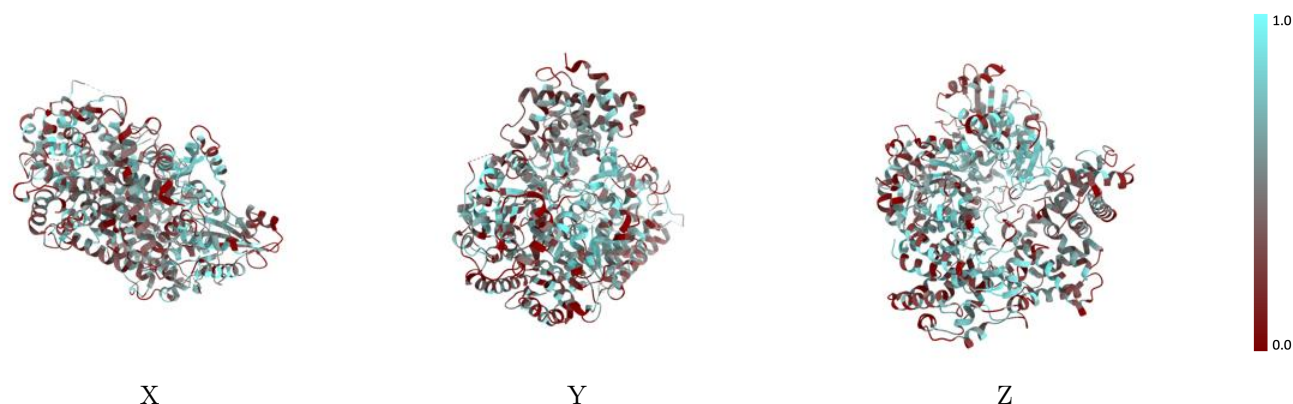
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



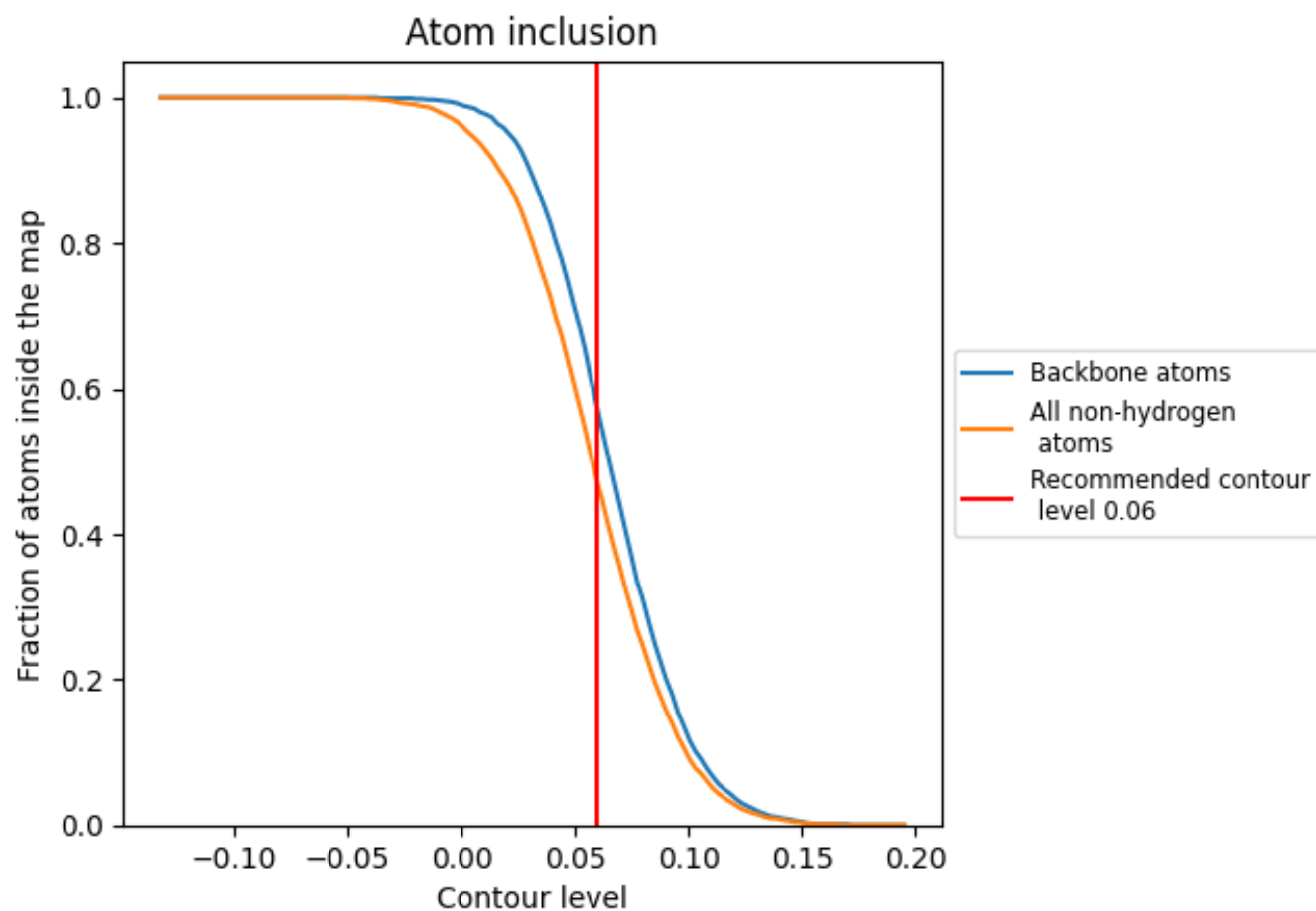
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).

9.4 Atom inclusion [i](#)



At the recommended contour level, 57% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4700	0.2800
A	0.5050	0.3190
B	0.4150	0.2090
C	0.3770	0.1930

