



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 06:37 AM UTC

PDB ID : 9O82 / pdb_00009o82
EMDB ID : EMD-70215
Title : Cryo-EM structure of apo rabbit TRPM3 having 2 resting and 2 activated subunits (para position) at 37 degrees Celsius
Authors : Kumar, S.; Lu, W.; Du, J.
Deposited on : 2025-04-15
Resolution : 4.46 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

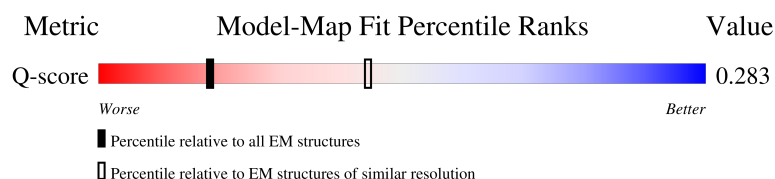
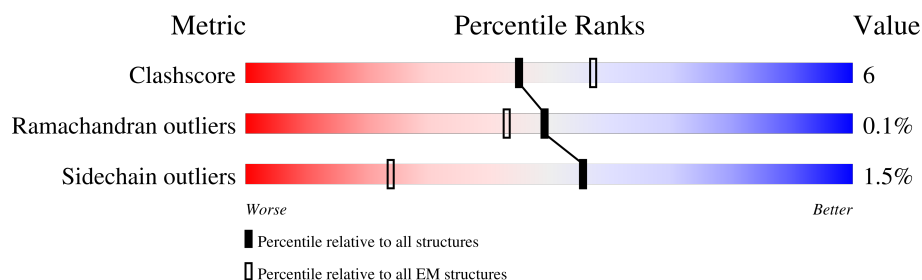
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.46 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	3001 (3.96 - 4.96)

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	1306	19% 65% 14% 21%
1	B	1306	24% 67% 11% 21%
1	C	1306	19% 66% 12% 21%
1	D	1306	24% 68% 11% 21%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 31311 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.

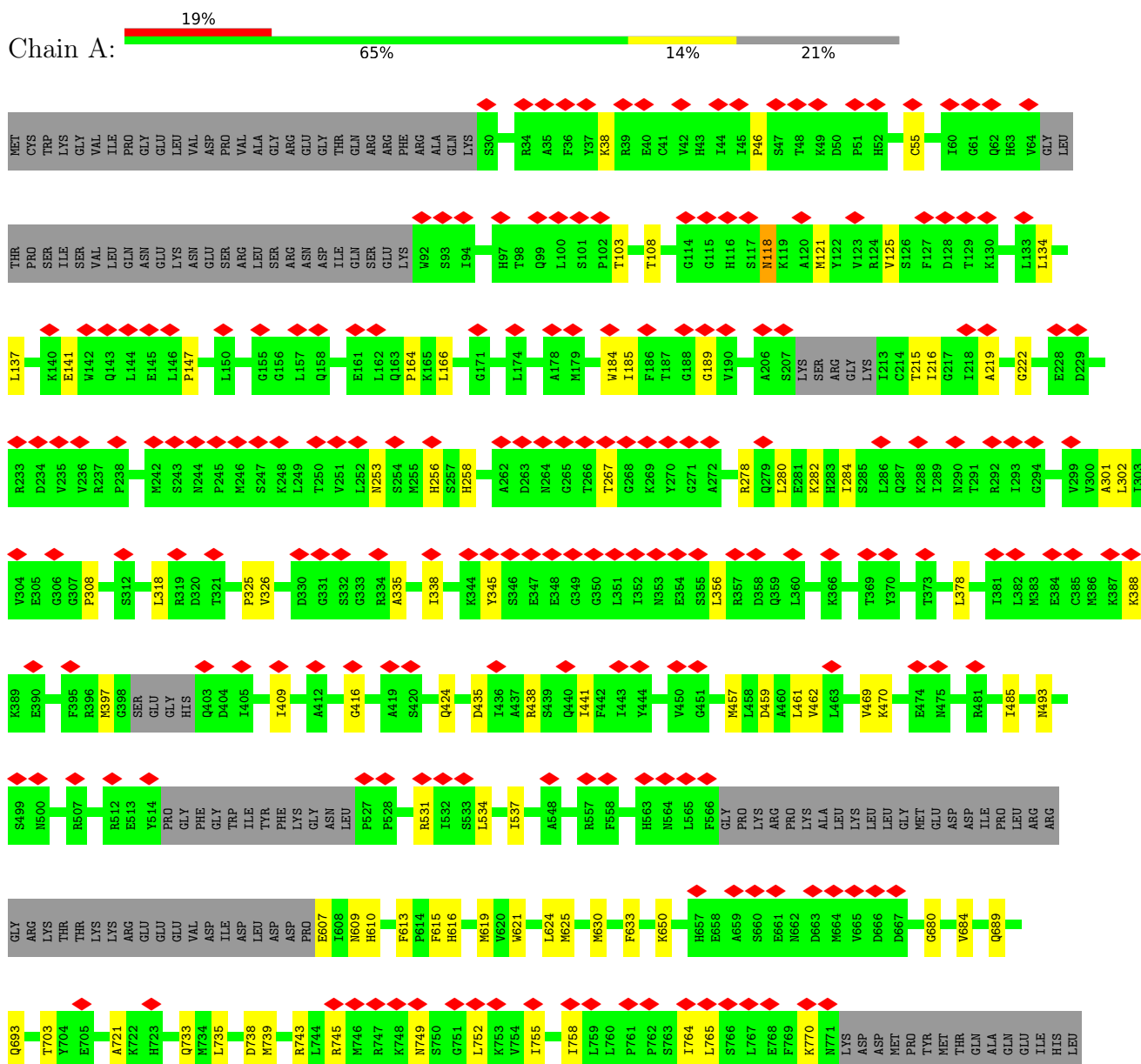
- Molecule 1 is a protein called TRPM3.

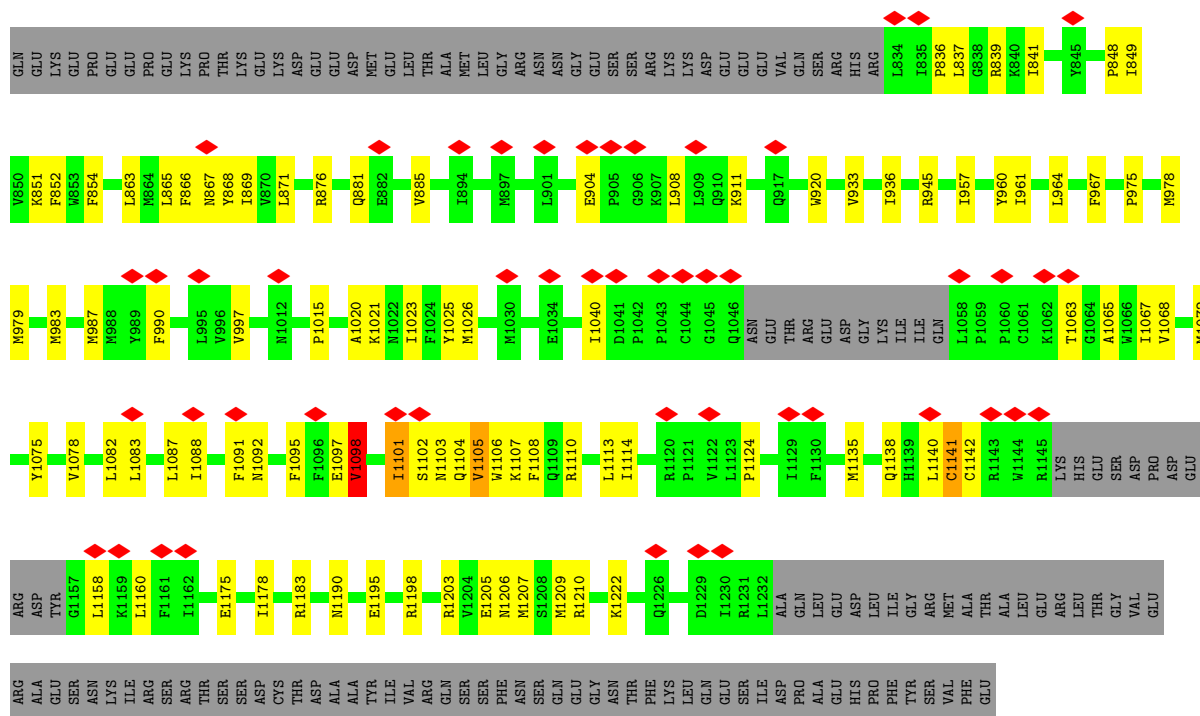
Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1031	Total	C	N	O	S	0	0
			7843	5078	1333	1379	53		
1	C	1030	Total	C	N	O	S	0	0
			7888	5109	1337	1385	57		
1	D	1033	Total	C	N	O	S	1	0
			7771	5043	1315	1357	56		
1	B	1036	Total	C	N	O	S	1	0
			7809	5061	1326	1365	57		

3 Residue-property plots

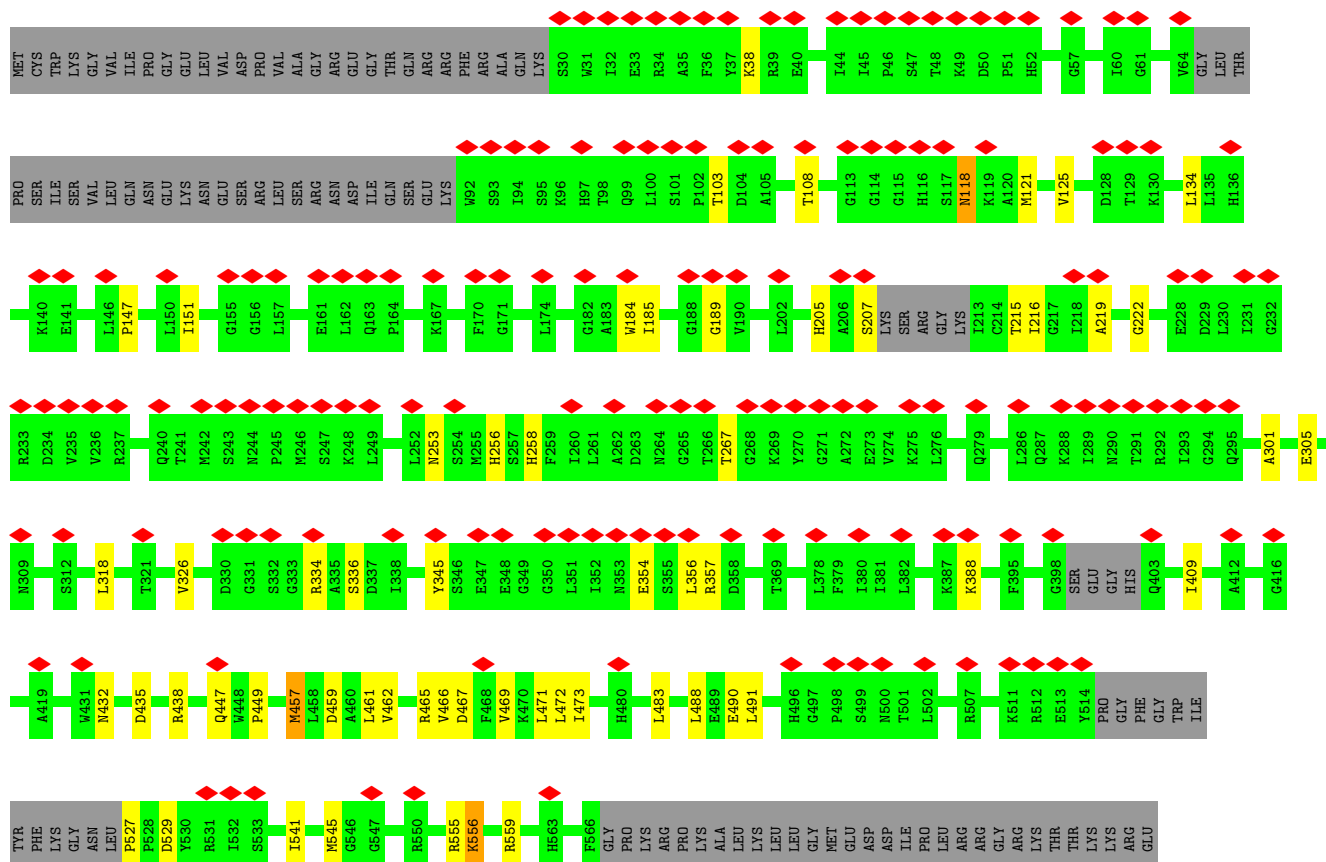
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

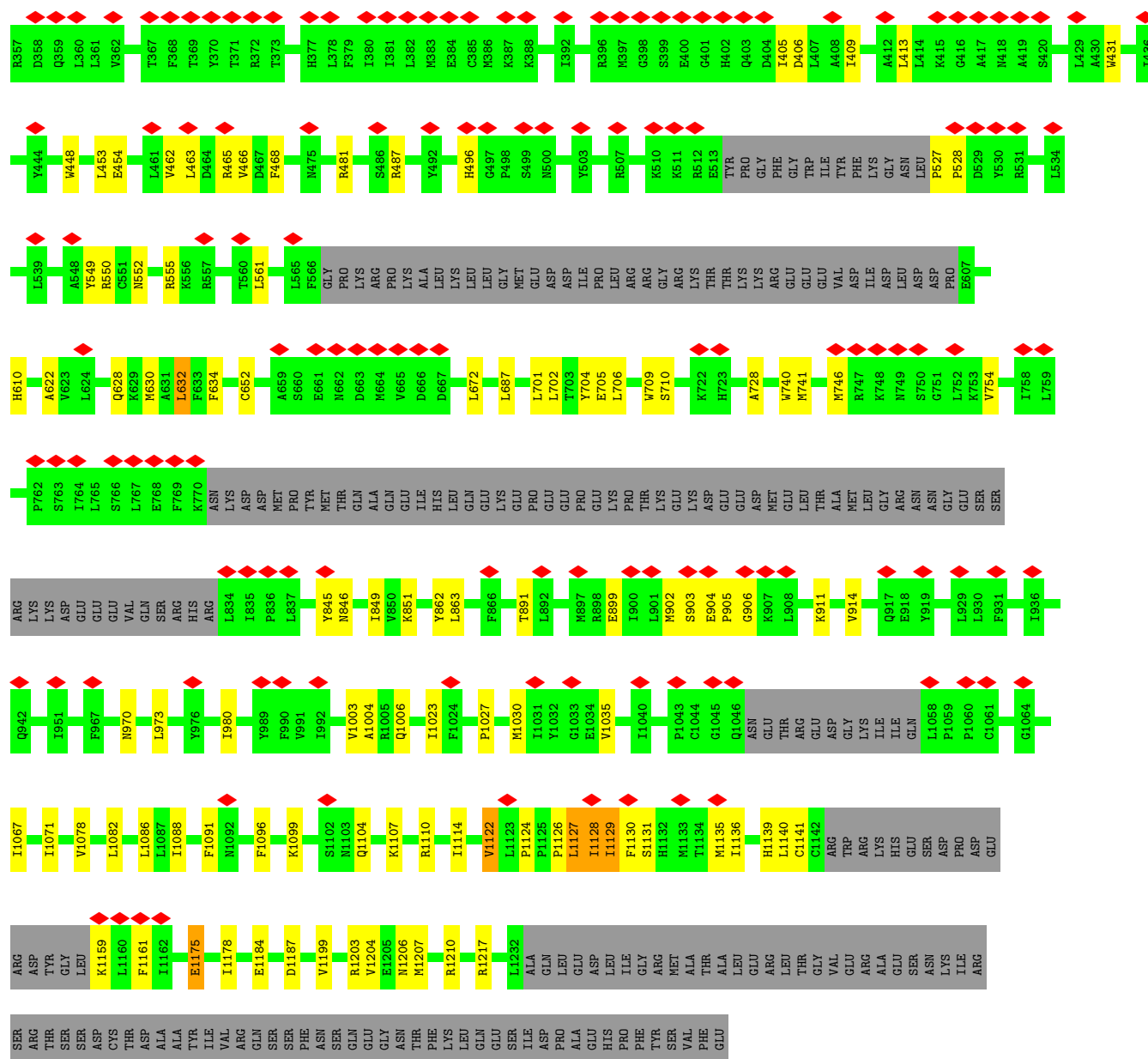
• Molecule 1: TRPM3



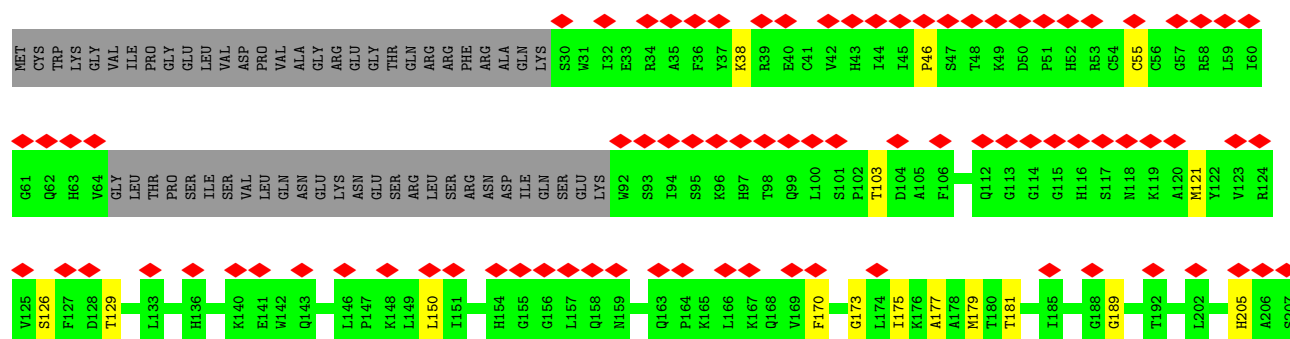


• Molecule 1: TRPM3





• Molecule 1: TRPM3





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	30726	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	58	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	1400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.008	Depositor
Minimum map value	-0.004	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.0024	Depositor
Map size ($\text{\AA}$)	317.184, 317.184, 317.184	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.826, 0.826, 0.826	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.42	3/8019 (0.0%)	0.86	6/10905 (0.1%)
1	B	0.39	2/7990 (0.0%)	0.80	10/10877 (0.1%)
1	C	0.42	1/8065 (0.0%)	0.85	8/10960 (0.1%)
1	D	0.41	0/7953	0.84	5/10830 (0.0%)
All	All	0.41	6/32027 (0.0%)	0.84	29/43572 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	904	GLU	C-N	6.55	1.41	1.33
1	A	1068	VAL	C-N	6.45	1.42	1.34
1	C	1026	MET	C-N	6.15	1.41	1.33
1	B	1026	MET	C-N	6.06	1.41	1.33
1	A	1026	MET	C-N	5.67	1.41	1.34
1	B	421	ALA	C-N	5.23	1.40	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1067	ILE	CA-C-N	8.86	125.82	120.24
1	B	1067	ILE	C-N-CA	8.86	125.82	120.24
1	D	1128	ILE	N-CA-C	-8.17	101.88	113.07
1	A	1098	VAL	N-CA-CB	7.74	119.61	110.55
1	A	868	TYR	CA-C-N	-7.42	113.02	121.85
1	A	868	TYR	C-N-CA	-7.42	113.02	121.85
1	D	1129	ILE	N-CA-C	-7.37	102.96	111.00
1	D	549	TYR	N-CA-C	-6.58	99.92	109.59
1	B	209	SER	CA-C-N	5.97	129.09	120.38
1	B	209	SER	C-N-CA	5.97	129.09	120.38
1	B	662	ASN	N-CA-C	-5.85	106.77	112.97
1	C	490	GLU	N-CA-CB	5.83	118.78	110.16
1	B	1093	ASN	CA-CB-CG	-5.78	106.82	112.60
1	C	556	LYS	CA-C-N	5.65	128.12	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	556	LYS	C-N-CA	5.65	128.12	120.38
1	B	1094	THR	N-CA-C	-5.59	105.67	112.88
1	A	738	ASP	N-CA-C	-5.57	104.90	110.97
1	C	1125	PRO	N-CA-C	5.50	117.41	110.70
1	A	979	MET	N-CA-C	-5.48	105.38	111.36
1	D	746	MET	CB-CA-C	-5.46	109.30	115.79
1	C	870	VAL	N-CA-C	-5.45	107.49	113.43
1	B	1024	PHE	CA-C-N	-5.44	114.40	122.56
1	B	1024	PHE	C-N-CA	-5.44	114.40	122.56
1	A	1098	VAL	CA-C-O	-5.23	115.63	121.17
1	B	1125	PRO	N-CA-C	5.22	117.07	110.70
1	D	1175	GLU	N-CA-CB	5.22	119.18	110.41
1	C	467	ASP	N-CA-C	5.04	117.15	111.11
1	C	457	MET	CB-CG-SD	-5.03	97.60	112.70
1	C	685	GLU	CA-CB-CG	5.02	124.14	114.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7843	0	7528	119	0
1	B	7809	0	7441	103	0
1	C	7888	0	7623	97	0
1	D	7771	0	7401	96	0
All	All	31311	0	29993	386	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 6.

All (386) 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:891:THR:CG2	1:D:1129:ILE:HD11	1.37	1.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:891:THR:HG22	1:D:1129:ILE:CD1	1.50	1.41
1:A:863:LEU:HD11	1:A:960:TYR:CE1	2.05	0.91
1:D:1124:PRO:HG2	1:D:1127:LEU:HB2	1.54	0.89
1:C:1086:LEU:HD21	1:D:980:ILE:HG22	1.61	0.82
1:A:1098:VAL:O	1:A:1102:SER:N	2.15	0.79
1:D:1128:ILE:HA	1:D:1131:SER:HB2	1.63	0.79
1:A:987:MET:HE2	1:D:1082:LEU:HB2	1.66	0.77
1:A:863:LEU:HD11	1:A:960:TYR:HE1	1.44	0.76
1:A:1021:LYS:O	1:A:1025:TYR:HB2	1.85	0.74
1:D:891:THR:HG21	1:D:1129:ILE:HD11	1.63	0.73
1:A:876:ARG:HD2	1:A:945:ARG:HH22	1.53	0.73
1:C:668:ILE:HA	1:C:671:GLU:HG3	1.72	0.70
1:D:208:LYS:O	1:D:208:LYS:HG3	1.90	0.70
1:A:1138:GLN:HA	1:A:1141:CYS:HB2	1.74	0.70
1:A:975:PRO:HB3	1:A:1102:SER:O	1.92	0.70
1:C:1197:ILE:HD11	1:B:1197:ILE:HD12	1.74	0.69
1:D:466:VAL:HG12	1:D:630:MET:HG3	1.75	0.69
1:D:1030:MET:HG3	1:D:1035:VAL:HA	1.76	0.68
1:B:208:LYS:HG3	1:B:208:LYS:O	1.92	0.68
1:C:960:TYR:HE1	1:B:996:VAL:HA	1.59	0.67
1:C:1219:HIS:HA	1:D:1217:ARG:HH22	1.57	0.67
1:A:863:LEU:HD11	1:A:960:TYR:CD1	2.29	0.67
1:A:1183:ARG:HH11	1:D:210:ARG:NH1	1.93	0.66
1:A:743:ARG:HD3	1:A:770:LYS:HB2	1.77	0.66
1:A:1183:ARG:HD2	1:D:210:ARG:NH1	2.11	0.66
1:D:1159:LYS:HD3	1:D:1161:PHE:H	1.61	0.65
1:A:854:PHE:HB3	1:A:1124:PRO:HG3	1.77	0.65
1:A:1203:ARG:HG2	1:A:1207:MET:HE2	1.78	0.65
1:C:1103:ASN:O	1:C:1107:LYS:HB3	1.96	0.65
1:A:1198:ARG:HG2	1:B:1196:ARG:HH12	1.61	0.64
1:D:891:THR:CG2	1:D:1129:ILE:CD1	2.33	0.64
1:D:846:ASN:HA	1:D:851:LYS:HE3	1.81	0.63
1:B:298:PRO:HB2	1:B:413:LEU:HD22	1.79	0.63
1:C:541:ILE:O	1:C:545:MET:HB3	2.00	0.62
1:D:891:THR:HG22	1:D:1129:ILE:HD11	0.65	0.62
1:C:466:VAL:HB	1:C:630:MET:HG2	1.83	0.61
1:B:938:ARG:HA	1:B:945:ARG:HA	1.82	0.61
1:C:1086:LEU:HD22	1:D:1091:PHE:HZ	1.66	0.61
1:B:655:MET:HG3	1:B:672:LEU:HD13	1.83	0.60
1:C:724:ARG:HA	1:C:1108:PHE:HD2	1.67	0.60
1:A:975:PRO:HB2	1:A:1102:SER:HB2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:691:TYR:HD2	1:C:731:CYS:HB2	1.66	0.60
1:D:38:LYS:HB3	1:D:103:THR:HG22	1.84	0.59
1:C:1103:ASN:O	1:C:1107:LYS:CB	2.51	0.59
1:D:298:PRO:HB2	1:D:413:LEU:HD22	1.84	0.59
1:C:838:GLY:O	1:C:842:TYR:HB2	2.02	0.59
1:D:181:THR:HG1	1:D:431:TRP:CD1	2.21	0.59
1:A:983:MET:HE1	1:A:1095:PHE:HA	1.85	0.59
1:B:457:MET:HE1	1:B:469:VAL:HG13	1.84	0.59
1:C:189:GLY:HA3	1:C:219:ALA:HB2	1.85	0.59
1:B:1122:VAL:HG13	1:B:1123:LEU:HG	1.84	0.59
1:C:1013:GLU:HG3	1:C:1019:LEU:HD21	1.84	0.58
1:A:216:ILE:HG13	1:A:258:HIS:HB2	1.85	0.58
1:B:441:ILE:HG13	1:B:448:TRP:CZ2	2.38	0.58
1:A:1110:ARG:HE	1:A:1114:ILE:HD11	1.67	0.58
1:D:1067:ILE:O	1:D:1071:ILE:HG13	2.04	0.58
1:A:457:MET:HE1	1:A:469:VAL:HG13	1.86	0.58
1:C:706:LEU:HB3	1:C:709:TRP:HB2	1.86	0.57
1:B:38:LYS:HB3	1:B:103:THR:HG22	1.86	0.57
1:B:464:ASP:HA	1:B:630:MET:CE	2.34	0.57
1:A:1203:ARG:O	1:A:1207:MET:HG2	2.05	0.57
1:B:436:ILE:O	1:B:440:GLN:N	2.37	0.57
1:C:216:ILE:HG13	1:C:258:HIS:HB2	1.86	0.57
1:D:970:ASN:HB3	1:D:973:LEU:HB2	1.86	0.57
1:A:764:ILE:HD13	1:A:841:ILE:HG22	1.85	0.57
1:C:659:ALA:HA	1:C:662:ASN:HB2	1.86	0.57
1:D:406:ASP:HA	1:D:409:ILE:HG22	1.86	0.57
1:A:1015:PRO:HD3	1:B:872:VAL:HG13	1.85	0.57
1:B:1005:ARG:HH22	1:B:1026:MET:HE2	1.70	0.57
1:C:665:VAL:HG23	1:C:668:ILE:HG12	1.87	0.56
1:A:613:PHE:HB3	1:A:616:HIS:HD1	1.70	0.56
1:C:1086:LEU:HD22	1:D:1091:PHE:CZ	2.40	0.56
1:C:723:HIS:CD2	1:C:726:PHE:HB2	2.41	0.56
1:B:882:GLU:HA	1:B:885:VAL:HG12	1.88	0.56
1:A:118:ASN:C	1:A:118:ASN:HD22	2.14	0.56
1:C:969:VAL:HG23	1:C:1113:LEU:HD21	1.88	0.55
1:A:1087:LEU:O	1:A:1091:PHE:HB2	2.06	0.55
1:A:189:GLY:HA3	1:A:219:ALA:HB2	1.87	0.55
1:A:534:LEU:HA	1:A:537:ILE:HG12	1.88	0.55
1:C:1021:LYS:O	1:C:1025:TYR:HB2	2.06	0.55
1:D:706:LEU:HB3	1:D:709:TRP:HB2	1.89	0.55
1:C:483:LEU:HD11	1:C:618:LEU:HD21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:LYS:O	1:D:208:LYS:CG	2.55	0.55
1:C:1101:ILE:HG23	1:C:1104:GLN:HB2	1.88	0.55
1:D:301:ALA:HB3	1:D:326:VAL:HG12	1.88	0.55
1:D:705:GLU:HB2	1:D:1159:LYS:HG3	1.87	0.55
1:A:493:ASN:HD22	1:A:531:ARG:HA	1.72	0.54
1:C:301:ALA:HB3	1:C:326:VAL:HG12	1.90	0.54
1:C:1221:MET:HE1	1:B:1222:LYS:HA	1.90	0.54
1:B:406:ASP:HA	1:B:409:ILE:HG22	1.88	0.54
1:A:1091:PHE:CE2	1:D:1086:LEU:HA	2.43	0.54
1:C:867:ASN:HA	1:C:870:VAL:HG12	1.88	0.54
1:D:189:GLY:HA3	1:D:219:ALA:HB2	1.88	0.54
1:A:1183:ARG:HB3	1:D:210:ARG:HH12	1.70	0.54
1:B:189:GLY:HA3	1:B:219:ALA:HB2	1.88	0.54
1:C:118:ASN:C	1:C:118:ASN:HD22	2.15	0.54
1:A:318:LEU:HD13	1:A:388:LYS:HB2	1.89	0.54
1:A:957:ILE:HD11	1:D:1004:ALA:HB2	1.90	0.54
1:A:1078:VAL:O	1:A:1082:LEU:HB3	2.07	0.54
1:C:459:ASP:HA	1:C:462:VAL:HG22	1.90	0.54
1:D:1107:LYS:HA	1:D:1110:ARG:HG3	1.89	0.54
1:D:1003:VAL:HG12	1:D:1023:ILE:HD12	1.89	0.54
1:B:208:LYS:O	1:B:208:LYS:CG	2.56	0.54
1:C:857:LEU:HA	1:C:860:ILE:HG22	1.89	0.53
1:B:441:ILE:HG23	1:B:448:TRP:HZ2	1.73	0.53
1:A:1082:LEU:HD21	1:B:980:ILE:HG23	1.90	0.53
1:A:865:LEU:HD12	1:A:881:GLN:HG3	1.89	0.53
1:C:432:ASN:HB3	1:C:465:ARG:HH12	1.73	0.53
1:A:863:LEU:CD1	1:A:960:TYR:HE1	2.19	0.53
1:C:851:LYS:HD3	1:C:1120:ARG:HB2	1.89	0.53
1:D:899:GLU:O	1:D:903:SER:HB3	2.09	0.53
1:A:621:TRP:CD1	1:A:625:MET:HE3	2.44	0.53
1:D:46:PRO:HG2	1:D:55:CYS:HB2	1.91	0.53
1:D:741:MET:HA	1:D:849:ILE:HD12	1.90	0.53
1:C:151:ILE:HG21	1:C:409:ILE:HD11	1.91	0.53
1:A:185:ILE:HB	1:A:215:THR:HA	1.90	0.52
1:C:354:GLU:HA	1:C:357:ARG:HG2	1.91	0.52
1:C:1095:PHE:HA	1:C:1098:VAL:HG22	1.91	0.52
1:D:300:VAL:HB	1:D:413:LEU:HD21	1.92	0.52
1:A:749:ASN:HB3	1:A:752:LEU:HD12	1.91	0.52
1:D:138:MET:HE1	1:D:216:ILE:HG12	1.91	0.52
1:D:150:LEU:HD21	1:D:281:GLU:HG2	1.91	0.52
1:B:301:ALA:HB3	1:B:326:VAL:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:607:GLU:HA	1:A:610:HIS:HD2	1.74	0.52
1:A:1082:LEU:HB2	1:B:987:MET:HE1	1.90	0.52
1:A:302:LEU:HD21	1:A:409:ILE:HD12	1.92	0.52
1:B:764:ILE:HG21	1:B:841:ILE:HG22	1.92	0.52
1:A:345:TYR:HB3	1:A:356:LEU:HD13	1.91	0.52
1:C:1121:PRO:HB2	1:C:1128:ILE:HD11	1.92	0.52
1:B:619:MET:HE2	1:B:644:LYS:HE2	1.91	0.52
1:B:46:PRO:HG2	1:B:55:CYS:HB2	1.91	0.51
1:C:185:ILE:HB	1:C:215:THR:HA	1.91	0.51
1:C:318:LEU:HD13	1:C:388:LYS:HB2	1.93	0.51
1:A:38:LYS:HD3	1:A:103:THR:HA	1.91	0.51
1:A:983:MET:CE	1:A:1098:VAL:HB	2.40	0.51
1:B:494:THR:HG22	1:B:496:HIS:H	1.75	0.51
1:C:125:VAL:HG11	1:C:134:LEU:HD21	1.93	0.51
1:B:697:LEU:HD23	1:B:700:LYS:HZ3	1.76	0.51
1:A:1205:GLU:HA	1:B:1203:ARG:HH12	1.74	0.51
1:C:689:GLN:O	1:C:693:GLN:HB2	2.11	0.51
1:C:435:ASP:HA	1:C:438:ARG:HG2	1.93	0.50
1:B:903:SER:HB3	1:B:911:LYS:HB2	1.94	0.50
1:A:435:ASP:HA	1:A:438:ARG:HB3	1.93	0.50
1:C:38:LYS:HD3	1:C:103:THR:HA	1.93	0.50
1:D:465:ARG:HB3	1:D:468:PHE:HB2	1.93	0.50
1:B:1112:GLN:HA	1:B:1115:MET:HE2	1.93	0.50
1:C:873:LYS:HZ3	1:B:1014:GLU:HG2	1.76	0.50
1:C:1096:PHE:HA	1:C:1099:LYS:HE3	1.93	0.50
1:B:1035:VAL:HB	1:B:1040:ILE:HD11	1.94	0.50
1:B:300:VAL:HB	1:B:413:LEU:HD21	1.93	0.50
1:C:739:MET:HE1	1:C:1158:LEU:HA	1.94	0.50
1:C:743:ARG:HG3	1:C:770:LYS:HD3	1.92	0.50
1:C:438:ARG:HB3	1:C:471:LEU:HD13	1.94	0.50
1:B:640:GLU:HB3	1:B:643:ALA:HB3	1.93	0.49
1:B:1095:PHE:HA	1:B:1098:VAL:HG12	1.93	0.49
1:A:222:GLY:HA3	1:A:267:THR:HA	1.94	0.49
1:D:222:GLY:HA3	1:D:267:THR:HA	1.93	0.49
1:B:878:PRO:HB2	1:B:883:TRP:CD1	2.47	0.49
1:A:983:MET:HE3	1:A:1098:VAL:HG21	1.94	0.49
1:C:345:TYR:HB3	1:C:356:LEU:HD13	1.95	0.49
1:C:1095:PHE:CD2	1:B:1093:ASN:CG	2.90	0.49
1:A:1020:ALA:HA	1:A:1023:ILE:HD12	1.94	0.49
1:C:527:PRO:HG2	1:C:529:ASP:HB2	1.94	0.49
1:B:441:ILE:HG13	1:B:448:TRP:CH2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:170:PHE:HA	1:D:405:ILE:HG13	1.95	0.49
1:B:484:THR:H	1:B:487:ARG:HB3	1.76	0.49
1:A:108:THR:HA	1:A:121:MET:HA	1.94	0.48
1:B:181:THR:HA	1:B:431:TRP:HD1	1.78	0.48
1:D:454:GLU:HB2	1:D:487:ARG:HH12	1.77	0.48
1:A:758:ILE:HG23	1:A:854:PHE:HE1	1.78	0.48
1:B:870:VAL:HG21	1:B:956:ILE:HD11	1.94	0.48
1:B:552:ASN:HA	1:B:555:ARG:HG2	1.95	0.48
1:A:755:ILE:HA	1:A:758:ILE:HD12	1.96	0.48
1:B:216:ILE:HG13	1:B:258:HIS:HB2	1.96	0.48
1:B:222:GLY:HA3	1:B:267:THR:HA	1.96	0.48
1:B:181:THR:HA	1:B:431:TRP:CD1	2.49	0.48
1:A:301:ALA:HB3	1:A:326:VAL:HG23	1.96	0.48
1:A:1065:ALA:C	1:A:1067:ILE:H	2.21	0.48
1:B:999:MET:HE2	1:B:999:MET:HB2	1.75	0.48
1:A:650:LYS:HG2	1:A:721:ALA:HA	1.95	0.48
1:D:706:LEU:HD12	1:D:709:TRP:HD1	1.78	0.47
1:A:424:GLN:HB3	1:A:441:ILE:HD12	1.96	0.47
1:A:1175:GLU:HA	1:A:1178:ILE:HG22	1.96	0.47
1:C:1001:PHE:HD2	1:C:1027:PRO:HB3	1.79	0.47
1:D:153:VAL:HG12	1:D:302:LEU:HB3	1.96	0.47
1:A:703:THR:HG23	1:A:735:LEU:HD11	1.95	0.47
1:C:937:LEU:HD12	1:C:944:PHE:HB3	1.96	0.47
1:A:459:ASP:HA	1:A:462:VAL:HG22	1.95	0.47
1:A:983:MET:HE3	1:A:1098:VAL:CG2	2.45	0.47
1:A:185:ILE:HG13	1:A:215:THR:HG22	1.97	0.47
1:A:876:ARG:HA	1:A:945:ARG:HH12	1.79	0.47
1:B:1044:CYS:HA	1:B:1058:LEU:HB3	1.96	0.47
1:A:485:ILE:H	1:A:609:ASN:HB2	1.80	0.47
1:A:1106:TRP:O	1:A:1107:LYS:C	2.55	0.47
1:C:1086:LEU:HD13	1:D:1091:PHE:CE1	2.49	0.47
1:B:662:ASN:O	1:B:663:ASP:C	2.58	0.47
1:B:723:HIS:HD2	1:B:726:PHE:HB2	1.80	0.47
1:D:1027:PRO:HA	1:D:1030:MET:HE2	1.97	0.47
1:B:1005:ARG:HH21	1:B:1027:PRO:HD3	1.80	0.47
1:A:990:PHE:HE1	1:A:1083:LEU:HB3	1.79	0.47
1:C:555:ARG:HA	1:C:555:ARG:HH11	1.80	0.46
1:A:164:PRO:HD2	1:A:1209:MET:HE1	1.98	0.46
1:A:836:PRO:HA	1:A:839:ARG:HD3	1.97	0.46
1:D:216:ILE:HG13	1:D:258:HIS:HB2	1.97	0.46
1:C:1110:ARG:O	1:C:1114:ILE:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1096:PHE:HA	1:D:1099:LYS:HE3	1.97	0.46
1:D:1126:PRO:O	1:D:1129:ILE:HG12	2.15	0.46
1:B:170:PHE:HA	1:B:405:ILE:HG13	1.97	0.46
1:B:463:LEU:O	1:B:464:ASP:C	2.57	0.46
1:A:920:TRP:HH2	1:A:961:ILE:HG22	1.80	0.46
1:A:964:LEU:O	1:A:967:PHE:HB2	2.15	0.46
1:D:1006:GLN:HG3	1:D:1023:ILE:HD13	1.98	0.46
1:A:978:MET:HE3	1:A:1105:VAL:HB	1.97	0.46
1:A:1206:ASN:O	1:A:1210:ARG:HB3	2.16	0.46
1:A:745:ARG:HB3	1:A:770:LYS:HG3	1.96	0.46
1:B:150:LEU:HB3	1:B:299:VAL:HG12	1.98	0.46
1:C:222:GLY:HA3	1:C:267:THR:HA	1.96	0.46
1:C:915:TRP:HH2	1:C:925:LEU:HB2	1.80	0.46
1:D:652:CYS:SG	1:D:672:LEU:O	2.74	0.46
1:A:733:GLN:HB2	1:A:1107:LYS:NZ	2.31	0.45
1:A:866:PHE:HD1	1:A:885:VAL:HG11	1.82	0.45
1:D:179:MET:HE2	1:D:211:GLY:HA3	1.98	0.45
1:A:1104:GLN:O	1:A:1105:VAL:C	2.58	0.45
1:C:305:GLU:HG3	1:C:334:ARG:HB2	1.98	0.45
1:D:448:TRP:CZ2	1:D:453:LEU:HG	2.51	0.45
1:A:613:PHE:HB3	1:A:616:HIS:ND1	2.31	0.45
1:C:457:MET:HE1	1:C:469:VAL:HG13	1.97	0.45
1:C:1086:LEU:HD13	1:D:1091:PHE:CZ	2.52	0.45
1:D:108:THR:HB	1:D:236:VAL:HG22	1.98	0.45
1:D:1078:VAL:O	1:D:1082:LEU:HB3	2.16	0.45
1:C:185:ILE:HG13	1:C:215:THR:HG22	1.98	0.45
1:C:253:ASN:HB3	1:C:256:HIS:CE1	2.51	0.45
1:C:620:VAL:O	1:C:624:LEU:HB2	2.17	0.45
1:D:705:GLU:HB3	1:D:1159:LYS:HE3	1.99	0.45
1:B:1110:ARG:O	1:B:1114:ILE:HG12	2.17	0.45
1:A:908:LEU:HA	1:A:911:LYS:HG2	1.99	0.45
1:B:484:THR:HG22	1:B:609:ASN:HB2	1.98	0.45
1:B:505:LEU:HD12	1:B:544:LEU:HD21	1.99	0.45
1:C:983:MET:HB3	1:C:1091:PHE:CE1	2.52	0.45
1:B:485:ILE:HD12	1:B:609:ASN:HA	1.99	0.45
1:D:1175:GLU:HA	1:D:1178:ILE:HG22	1.99	0.45
1:D:845:TYR:HE2	1:D:1122:VAL:HG22	1.82	0.45
1:C:1002:GLY:HA3	1:C:1023:ILE:HG23	1.99	0.44
1:D:1159:LYS:NZ	1:D:1161:PHE:HA	2.32	0.44
1:B:173:GLY:HA3	1:B:405:ILE:HB	1.98	0.44
1:A:689:GLN:O	1:A:693:GLN:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:VAL:HG11	1:A:134:LEU:HD21	1.98	0.44
1:A:470:LYS:HG2	1:A:633:PHE:CE1	2.53	0.44
1:A:1222:LYS:HD3	1:A:1222:LYS:HA	1.68	0.44
1:D:1184:GLU:HA	1:D:1187:ASP:HB2	1.99	0.44
1:B:441:ILE:HG23	1:B:448:TRP:CZ2	2.52	0.44
1:B:1166:GLU:HA	1:B:1169:LYS:HG2	1.99	0.44
1:D:622:ALA:HB2	1:D:634:PHE:CD2	2.53	0.44
1:A:624:LEU:HD22	1:A:625:MET:HE2	2.00	0.44
1:C:873:LYS:NZ	1:B:1014:GLU:HG2	2.32	0.44
1:D:462:VAL:HG23	1:D:463:LEU:HG	1.99	0.44
1:A:863:LEU:CD1	1:A:960:TYR:CE1	2.90	0.44
1:A:983:MET:HE1	1:A:1098:VAL:HB	1.98	0.44
1:A:1190:ASN:HA	1:A:1195:GLU:HG2	1.99	0.44
1:C:447:GLN:HG3	1:C:449:PRO:HD3	2.00	0.44
1:C:723:HIS:HD2	1:C:726:PHE:HB2	1.80	0.44
1:C:883:TRP:HA	1:C:886:ILE:HG22	2.00	0.44
1:D:175:ILE:HD12	1:D:205:HIS:HB2	2.00	0.44
1:D:906:GLY:O	1:D:911:LYS:NZ	2.50	0.44
1:A:733:GLN:HB2	1:A:1107:LYS:HZ1	1.83	0.44
1:C:872:VAL:HG13	1:B:1015:PRO:HD3	1.98	0.44
1:D:126:SER:HB3	1:D:129:THR:HG23	1.99	0.44
1:C:108:THR:HA	1:C:121:MET:HA	2.00	0.44
1:B:694:ASP:HB3	1:B:697:LEU:HB2	2.00	0.44
1:B:177:ALA:O	1:B:181:THR:HG22	2.18	0.43
1:C:147:PRO:HB3	1:C:184:TRP:CD1	2.52	0.43
1:D:1204:VAL:HA	1:D:1207:MET:HG3	1.99	0.43
1:C:989:TYR:HA	1:C:992:ILE:HD12	2.01	0.43
1:C:1003:VAL:HG12	1:C:1023:ILE:HG12	2.01	0.43
1:B:863:LEU:HD11	1:B:960:TYR:HD1	1.84	0.43
1:A:739:MET:HG3	1:A:1158:LEU:HD21	2.00	0.43
1:A:866:PHE:HA	1:A:869:ILE:HG22	2.01	0.43
1:A:1183:ARG:HH11	1:D:210:ARG:HH11	1.63	0.43
1:B:421:ALA:HA	1:B:441:ILE:HD11	2.00	0.43
1:A:253:ASN:HB3	1:A:256:HIS:CE1	2.53	0.43
1:A:1065:ALA:C	1:A:1067:ILE:N	2.76	0.43
1:B:410:LEU:HA	1:B:413:LEU:HD12	2.01	0.43
1:C:1123:LEU:HD23	1:C:1123:LEU:HA	1.84	0.43
1:D:552:ASN:HA	1:D:555:ARG:HG2	2.00	0.43
1:B:854:PHE:HE2	1:B:1124:PRO:HD3	1.83	0.43
1:A:867:ASN:O	1:A:871:LEU:CB	2.66	0.43
1:B:440:GLN:HB3	1:B:441:ILE:H	1.73	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:869:ILE:HG21	1:C:882:GLU:HB2	2.00	0.43
1:D:740:TRP:CE3	1:D:1114:ILE:HG12	2.54	0.43
1:B:984:MET:HE3	1:B:984:MET:HB3	1.94	0.43
1:B:997:VAL:HG22	1:B:1075:TYR:HD1	1.84	0.43
1:D:527:PRO:HA	1:D:528:PRO:HD3	1.91	0.42
1:B:713:THR:HG22	1:B:715:LEU:H	1.83	0.42
1:A:1106:TRP:C	1:A:1108:PHE:N	2.77	0.42
1:C:933:VAL:HA	1:C:936:ILE:HG22	2.01	0.42
1:B:436:ILE:O	1:B:437:ALA:C	2.62	0.42
1:B:706:LEU:HB3	1:B:709:TRP:HB2	2.01	0.42
1:A:933:VAL:HA	1:A:936:ILE:HG22	2.01	0.42
1:D:152:SER:HB3	1:D:277:ARG:HH21	1.84	0.42
1:D:754:VAL:HG11	1:D:849:ILE:HG22	2.01	0.42
1:B:282:LYS:HD2	1:B:282:LYS:HA	1.78	0.42
1:B:438:ARG:HA	1:B:442:PHE:HB2	2.00	0.42
1:B:746:MET:HE3	1:B:746:MET:HB3	1.62	0.42
1:C:613:PHE:HB3	1:C:616:HIS:HD1	1.83	0.42
1:C:1078:VAL:HG23	1:C:1082:LEU:HD12	2.01	0.42
1:B:761:PRO:N	1:B:762:PRO:HD2	2.34	0.42
1:D:481:ARG:HA	1:D:610:HIS:CE1	2.54	0.42
1:D:622:ALA:HB2	1:D:634:PHE:HD2	1.85	0.42
1:B:687:LEU:HA	1:B:690:SER:HB2	2.01	0.42
1:D:462:VAL:HB	1:D:496:HIS:HB2	2.01	0.42
1:D:1199:VAL:HG22	1:D:1203:ARG:HE	1.84	0.42
1:A:278:ARG:HG3	1:A:282:LYS:HE2	2.02	0.42
1:A:325:PRO:HG3	1:A:416:GLY:HA3	2.01	0.42
1:D:550:ARG:HH11	1:D:555:ARG:NH2	2.18	0.42
1:B:1139:HIS:HA	1:B:1142:CYS:HB2	2.02	0.42
1:A:147:PRO:HB3	1:A:184:TRP:CD1	2.55	0.42
1:A:308:PRO:HB3	1:A:378:LEU:HD21	2.02	0.42
1:B:969:VAL:HG13	1:B:1113:LEU:HD21	2.01	0.42
1:C:1183:ARG:HB3	1:B:210:ARG:HH12	1.84	0.42
1:B:150:LEU:HD21	1:B:281:GLU:HG2	2.01	0.42
1:A:335:ALA:HA	1:A:338:ILE:HG12	2.00	0.42
1:A:1040:ILE:HD12	1:A:1072:MET:HE2	2.02	0.42
1:A:1088:ILE:O	1:A:1092:ASN:ND2	2.53	0.42
1:A:1135:MET:HE2	1:A:1135:MET:HB3	1.96	0.42
1:D:701:LEU:HD12	1:D:704:TYR:HE1	1.85	0.42
1:A:137:LEU:HD12	1:A:141:GLU:HB2	2.01	0.41
1:C:205:HIS:CE1	1:C:207:SER:HB2	2.55	0.41
1:C:740:TRP:HH2	1:C:852:PHE:HB2	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1019:LEU:O	1:C:1023:ILE:HD12	2.21	0.41
1:D:905:PRO:HD3	1:D:914:VAL:HG21	2.02	0.41
1:B:436:ILE:HB	1:B:440:GLN:HB2	2.01	0.41
1:B:458:LEU:HD13	1:B:487:ARG:HD2	2.02	0.41
1:B:697:LEU:HD23	1:B:697:LEU:HA	1.89	0.41
1:B:866:PHE:HB2	1:B:885:VAL:HG21	2.02	0.41
1:A:615:PHE:O	1:A:619:MET:HB2	2.20	0.41
1:D:550:ARG:H	1:D:710:SER:HB2	1.85	0.41
1:B:748:LYS:O	1:B:749:ASN:C	2.63	0.41
1:A:46:PRO:HG2	1:A:55:CYS:HB2	2.02	0.41
1:D:628:GLN:HG3	1:D:632:LEU:HD22	2.03	0.41
1:B:442:PHE:HA	1:B:448:TRP:HE1	1.84	0.41
1:B:724:ARG:HD2	1:B:1105:VAL:HG12	2.01	0.41
1:A:461:LEU:HA	1:A:630:MET:HE1	2.02	0.41
1:A:849:ILE:HG13	1:A:852:PHE:HE1	1.86	0.41
1:C:305:GLU:H	1:C:336:SER:HB3	1.86	0.41
1:D:327:VAL:HG21	1:D:409:ILE:HA	2.03	0.41
1:B:536:ASP:HA	1:B:539:LEU:HD12	2.00	0.41
1:B:639:GLU:HG3	1:B:1171:HIS:ND1	2.36	0.41
1:A:765:LEU:HD11	1:A:837:LEU:HD11	2.01	0.41
1:C:1082:LEU:HA	1:C:1082:LEU:HD23	1.84	0.41
1:D:687:LEU:HD11	1:D:702:LEU:HD11	2.03	0.41
1:A:1087:LEU:HD13	1:A:1087:LEU:HA	1.88	0.41
1:C:740:TRP:CD2	1:C:1114:ILE:HD12	2.56	0.41
1:D:728:ALA:HB2	1:D:1104:GLN:HG3	2.02	0.41
1:A:848:PRO:HA	1:A:851:LYS:HB2	2.01	0.41
1:A:1088:ILE:HG21	1:B:1088:ILE:HG23	2.02	0.41
1:C:668:ILE:HB	1:C:672:LEU:HD23	2.03	0.41
1:C:976:TYR:O	1:C:980:ILE:HG12	2.21	0.41
1:B:179:MET:SD	1:B:211:GLY:HA3	2.61	0.41
1:A:1097:GLU:OE2	1:A:1101:ILE:HG13	2.21	0.41
1:B:121:MET:HB2	1:B:258:HIS:CD2	2.56	0.41
1:B:175:ILE:HD12	1:B:205:HIS:HB2	2.03	0.41
1:B:427:LEU:HD13	1:B:431:TRP:HZ3	1.86	0.41
1:A:166:LEU:HD22	1:A:397:MET:HE2	2.02	0.40
1:C:1001:PHE:CD2	1:C:1027:PRO:HB3	2.57	0.40
1:D:973:LEU:HD23	1:D:973:LEU:HA	1.94	0.40
1:B:414:LEU:HG	1:B:440:GLN:HG3	2.03	0.40
1:A:1183:ARG:HD2	1:D:210:ARG:CZ	2.50	0.40
1:C:461:LEU:HD12	1:C:491:LEU:HD22	2.02	0.40
1:C:930:LEU:HB3	1:C:955:ASN:HD21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:146:LEU:HD23	1:D:146:LEU:HA	1.92	0.40
1:D:1128:ILE:HA	1:D:1131:SER:CB	2.44	0.40
1:D:1206:ASN:O	1:D:1210:ARG:HG2	2.20	0.40
1:A:997:VAL:HG22	1:A:1075:TYR:HD1	1.86	0.40
1:C:473:ILE:HD13	1:C:473:ILE:HA	1.91	0.40
1:C:1001:PHE:HD1	1:C:1071:ILE:HG22	1.87	0.40
1:D:862:TYR:HD2	1:D:863:LEU:HD22	1.85	0.40
1:B:1193:ASN:O	1:B:1197:ILE:HG12	2.21	0.40
1:A:680:GLY:O	1:A:684:VAL:HG23	2.22	0.40
1:D:185:ILE:HB	1:D:215:THR:HA	2.03	0.40
1:D:706:LEU:HD12	1:D:709:TRP:CD1	2.57	0.40
1:B:126:SER:HB3	1:B:129:THR:HG23	2.02	0.40
1:A:280:LEU:O	1:A:284:ILE:HG12	2.21	0.40
1:C:556:LYS:HA	1:C:559:ARG:HG2	2.02	0.40
1:C:980:ILE:HG23	1:B:1082:LEU:HD21	2.04	0.40
1:C:1087:LEU:HD12	1:C:1087:LEU:HA	1.81	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	1013/1306 (78%)	950 (94%)	63 (6%)	0	100	100
1	B	1023/1306 (78%)	962 (94%)	60 (6%)	1 (0%)	48	83
1	C	1012/1306 (78%)	951 (94%)	60 (6%)	1 (0%)	48	83
1	D	1020/1306 (78%)	974 (96%)	45 (4%)	1 (0%)	48	83
All	All	4068/5224 (78%)	3837 (94%)	228 (6%)	3 (0%)	49	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	904	GLU
1	C	724	ARG
1	B	464	ASP

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	783/1162 (67%)	772 (99%)	11 (1%)	59	71
1	B	769/1162 (66%)	757 (98%)	12 (2%)	55	69
1	C	797/1162 (69%)	785 (98%)	12 (2%)	57	70
1	D	764/1162 (66%)	752 (98%)	12 (2%)	55	69
All	All	3113/4648 (67%)	3066 (98%)	47 (2%)	55	70

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	A	1063	THR
1	A	1098	VAL
1	A	1101	ILE
1	A	1103	ASN
1	A	1105	VAL
1	A	1113	LEU
1	A	1140	LEU
1	A	1141	CYS
1	A	1142	CYS
1	A	1160	LEU
1	C	118	ASN
1	C	472	LEU
1	C	488	LEU
1	C	863	LEU
1	C	965	ASP
1	C	1040	ILE
1	C	1086	LEU
1	C	1141	CYS

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Mol	Chain	Res	Type
1	C	1142	CYS
1	C	1176	GLN
1	C	1211	LEU
1	C	1213	GLU
1	D	561	LEU
1	D	632	LEU
1	D	902	MET
1	D	1088	ILE
1	D	1122	VAL
1	D	1127	LEU
1	D	1130	PHE
1	D	1135	MET
1	D	1136	ILE
1	D	1139	HIS
1	D	1140	LEU
1	D	1141	CYS
1	B	436	ILE
1	B	440	GLN
1	B	441	ILE
1	B	462	VAL
1	B	463	LEU
1	B	464	ASP
1	B	466	VAL
1	B	467	ASP
1	B	746	MET
1	B	999	MET
1	B	1086	LEU
1	B	1093	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	197	HIS
1	A	609	ASN
1	A	628	GLN
1	A	670	GLN
1	A	711	ASN
1	A	733	GLN
1	A	1046	GLN
1	A	1092	ASN
1	A	1215	ASN
1	C	118	ASN

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Mol	Chain	Res	Type
1	C	440	GLN
1	C	637	HIS
1	C	670	GLN
1	C	689	GLN
1	C	693	GLN
1	C	955	ASN
1	C	1092	ASN
1	C	1109	GLN
1	D	240	GLN
1	D	402	HIS
1	D	493	ASN
1	D	733	GLN
1	D	749	ASN
1	D	940	GLN
1	D	1109	GLN
1	D	1193	ASN
1	B	240	GLN
1	B	353	ASN
1	B	424	GLN
1	B	955	ASN
1	B	1022	ASN
1	B	1046	GLN
1	B	1215	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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

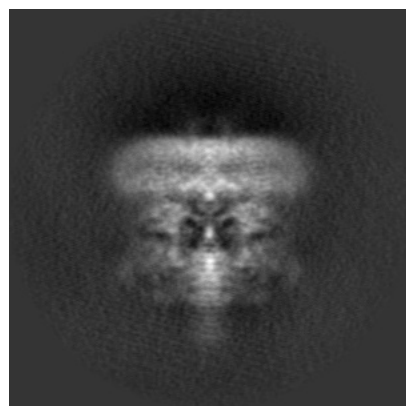
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70215. These allow visual inspection of the internal detail of the map and identification of artifacts.

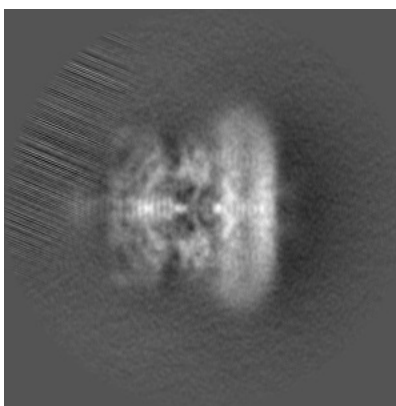
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

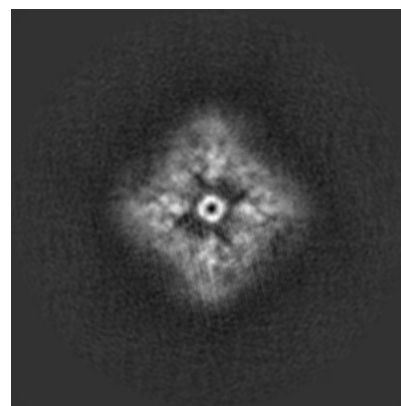
6.1.1 Primary map



X

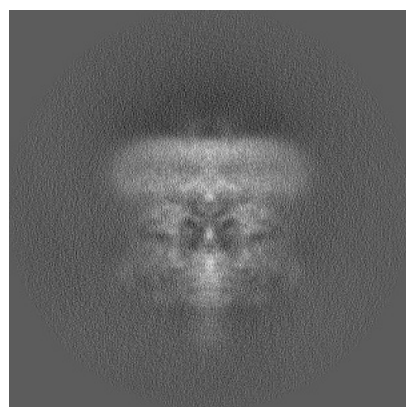


Y

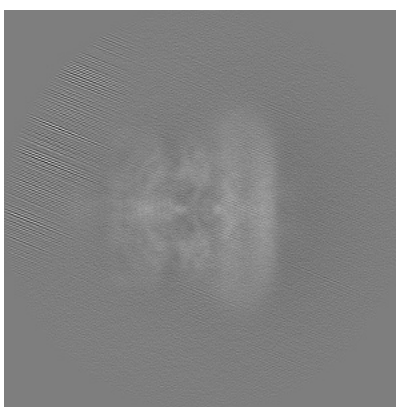


Z

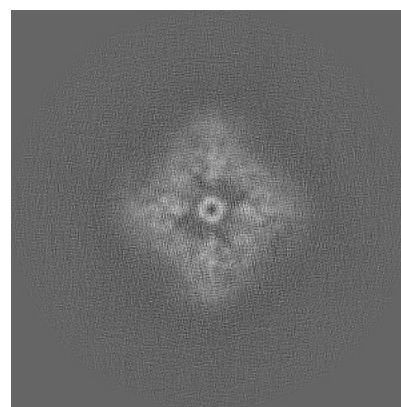
6.1.2 Raw map



X



Y

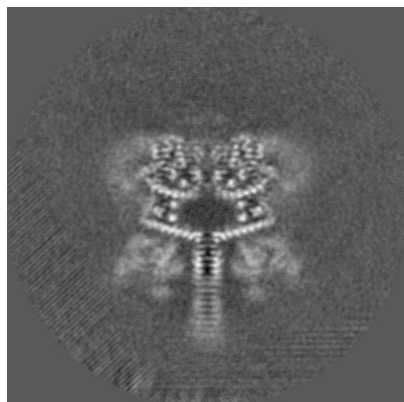


Z

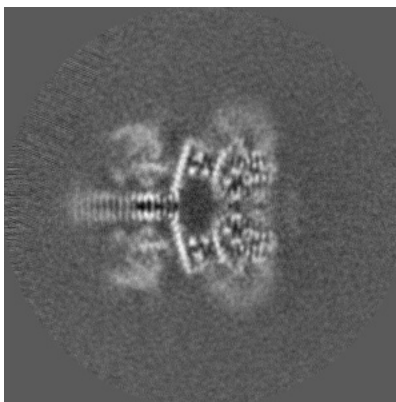
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

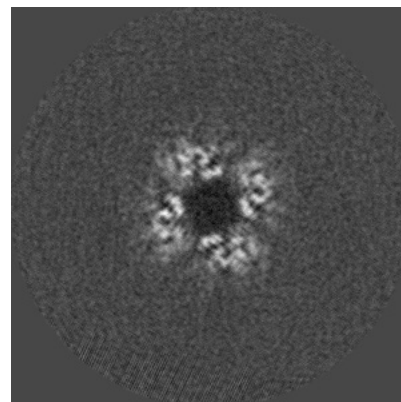
6.2.1 Primary map



X Index: 192

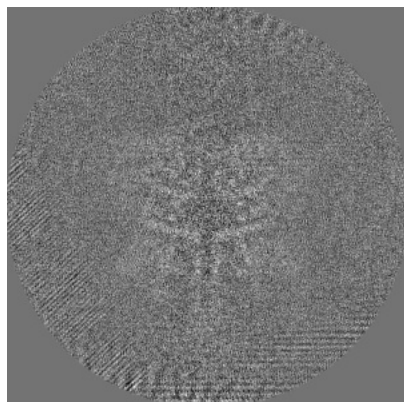


Y Index: 192

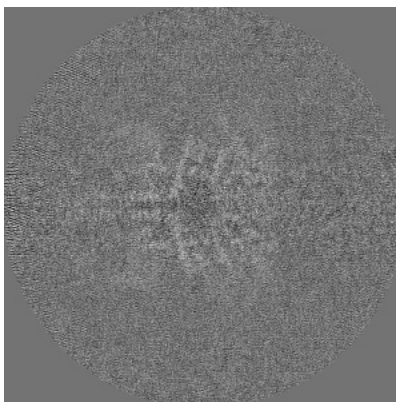


Z Index: 192

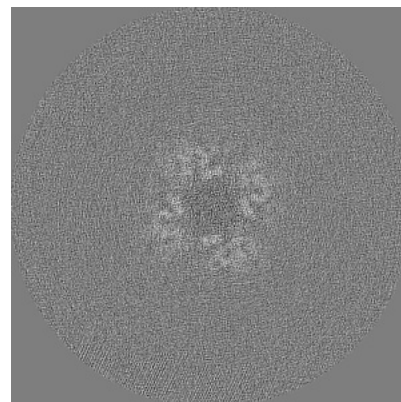
6.2.2 Raw map



X Index: 192



Y Index: 192

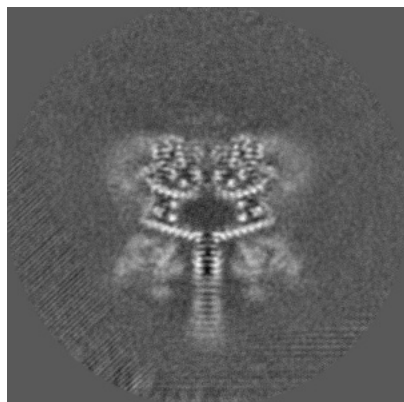


Z Index: 192

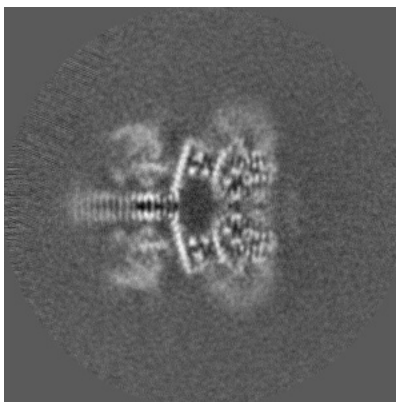
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

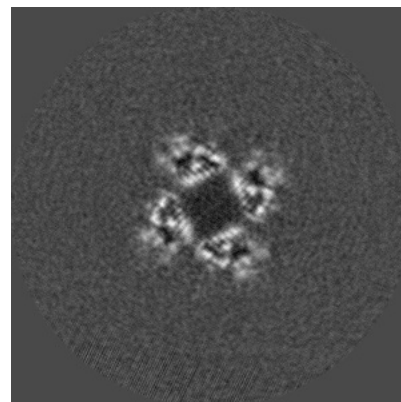
6.3.1 Primary map



X Index: 192

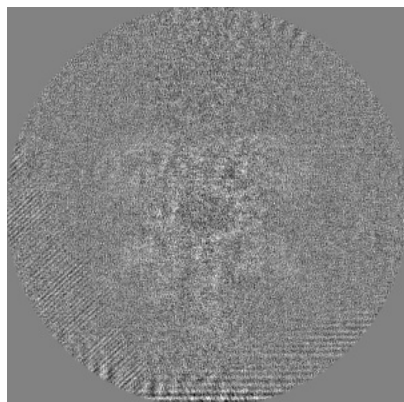


Y Index: 192

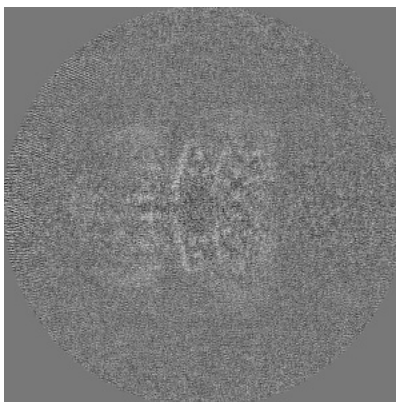


Z Index: 185

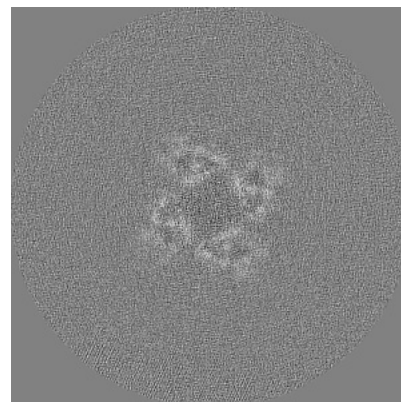
6.3.2 Raw map



X Index: 188



Y Index: 193

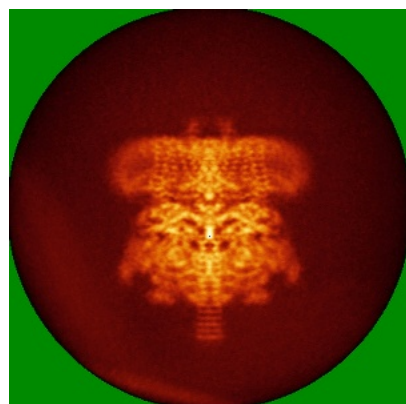


Z Index: 185

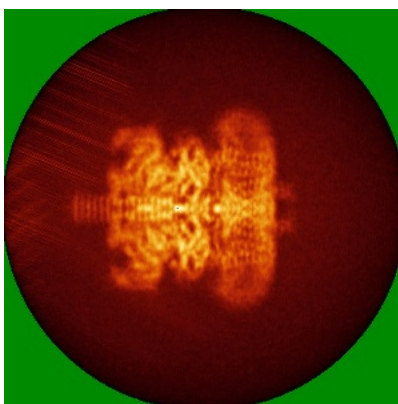
The images above show the largest variance slices of the map in three orthogonal directions.

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

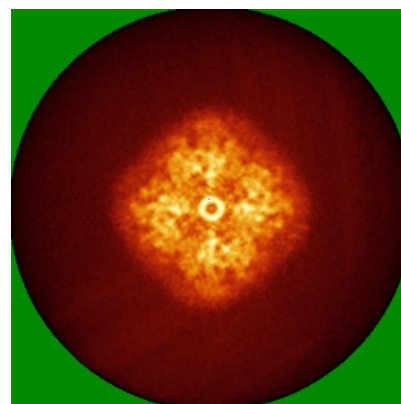
6.4.1 Primary map



X

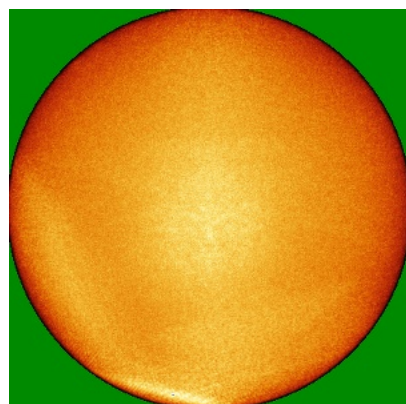


Y

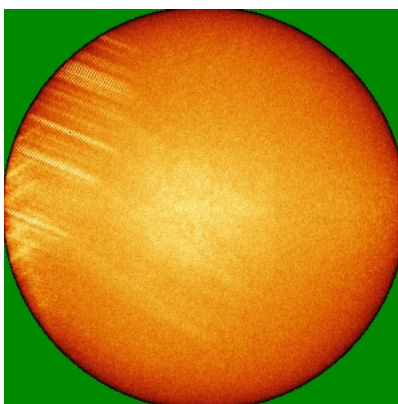


Z

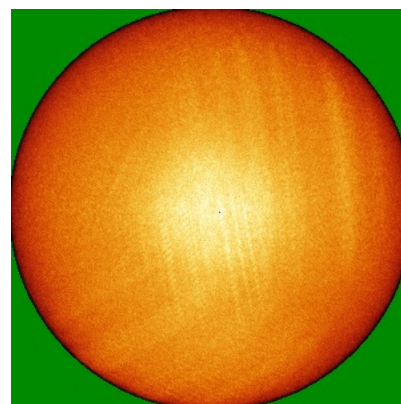
6.4.2 Raw map



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



X



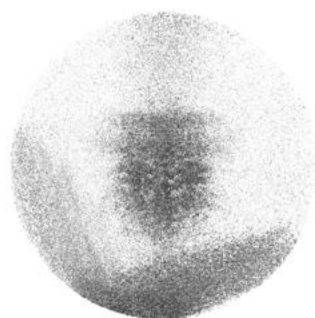
Y



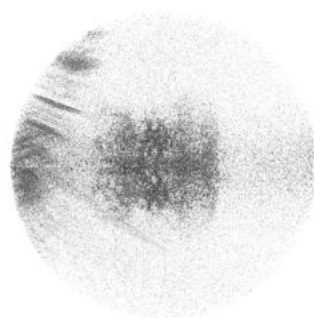
Z

The images above show the 3D surface view of the map at the recommended contour level 0.0024. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

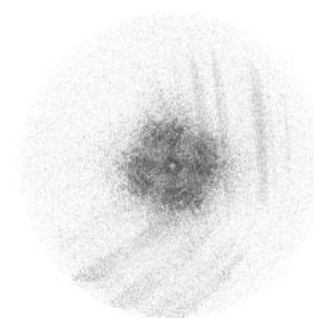
6.5.2 Raw map



X



Y



Z

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

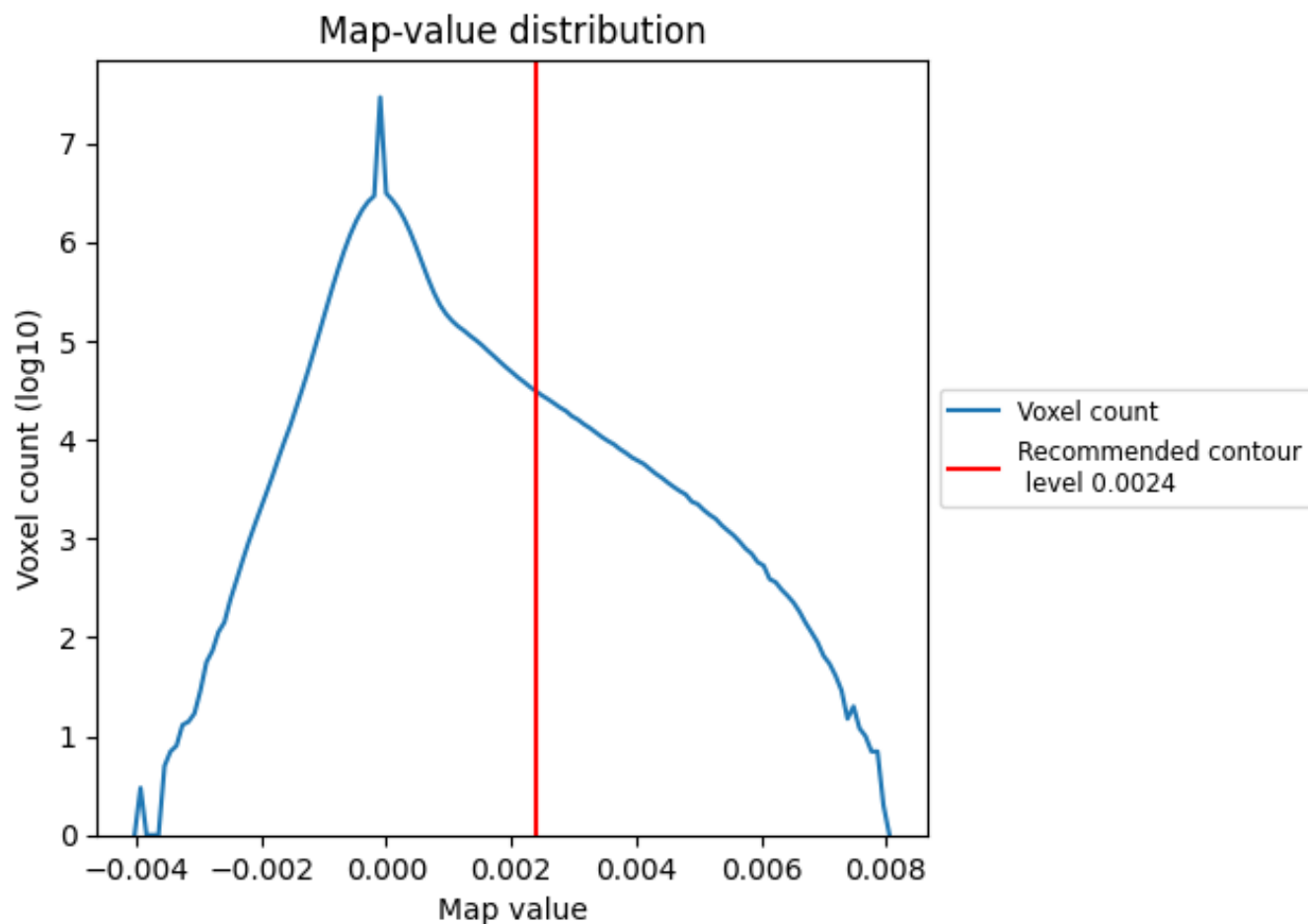
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

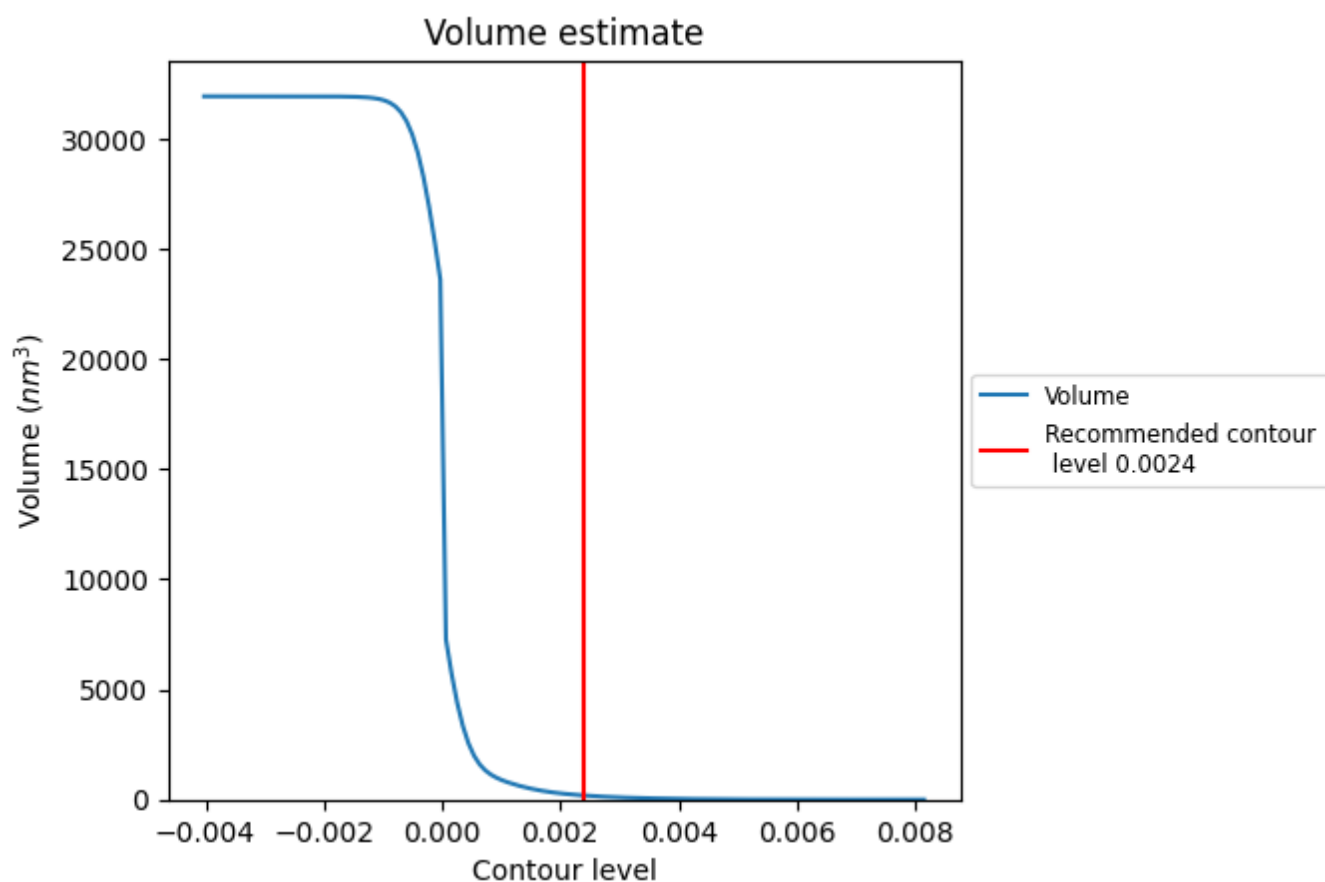
This section contains the results of statistical analysis of the map.

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.

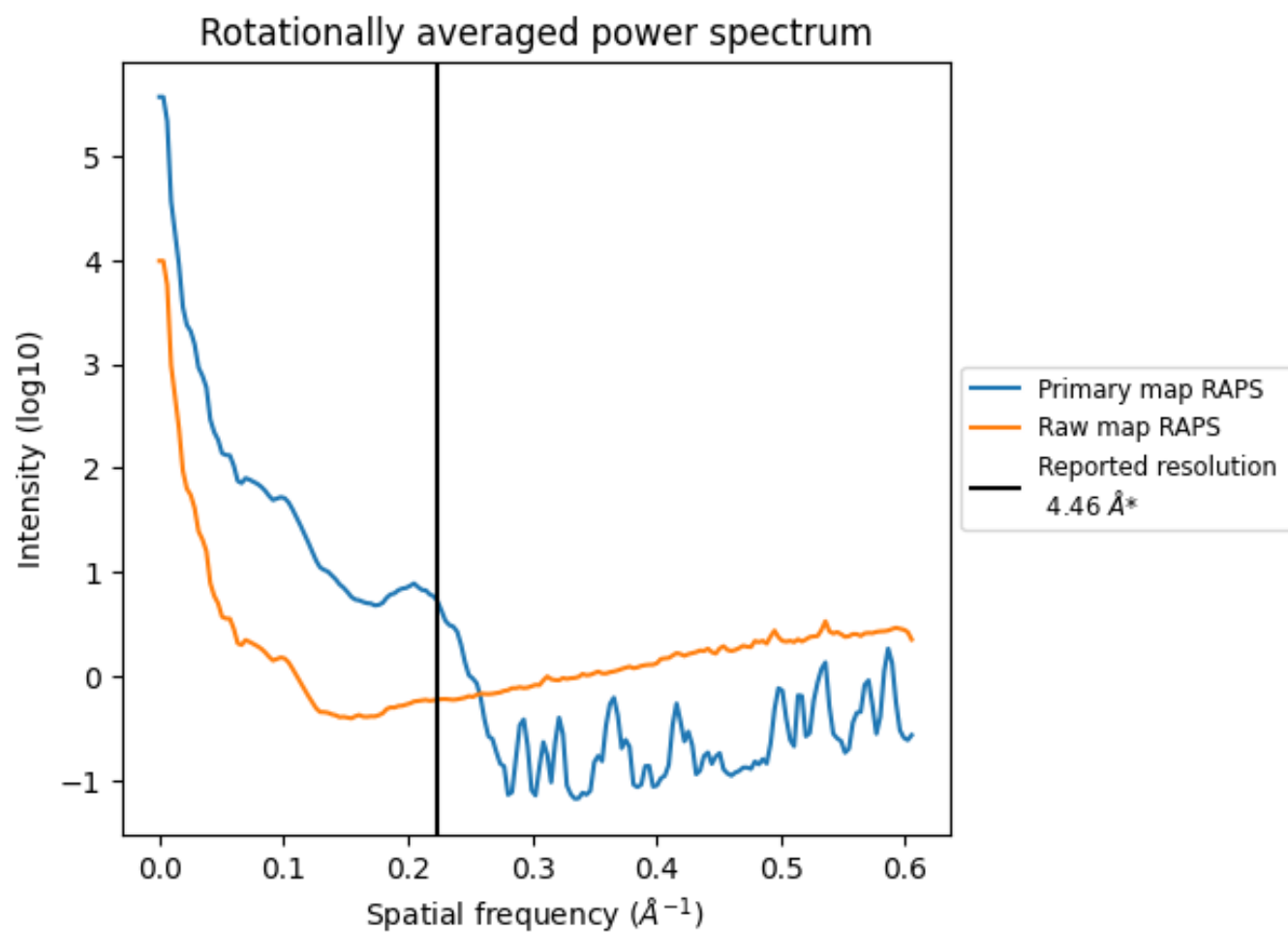
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 186 nm³; this corresponds to an approximate mass of 168 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 ⓘ

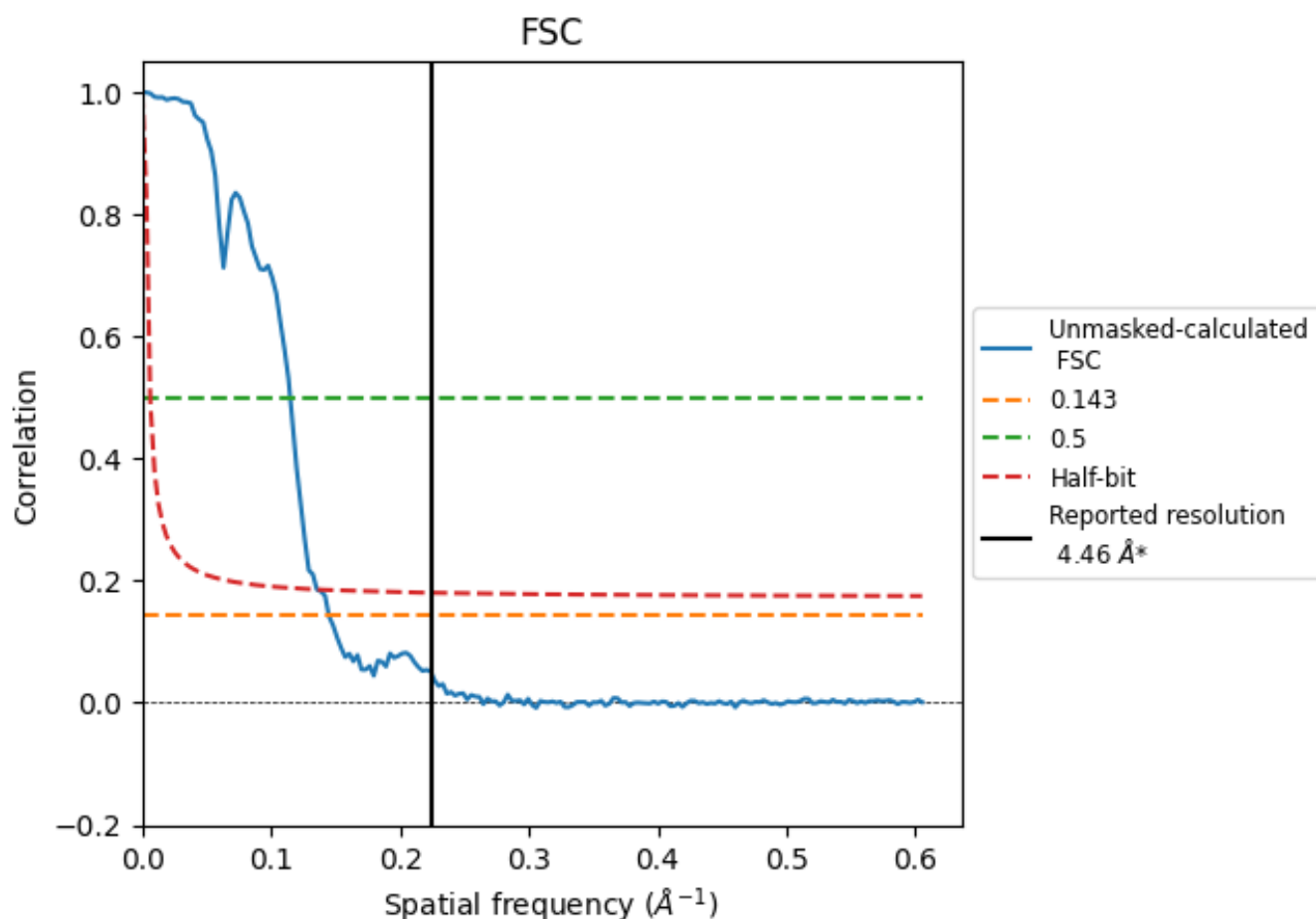


*Reported resolution corresponds to spatial frequency of 0.224 Å⁻¹

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 $0.224 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

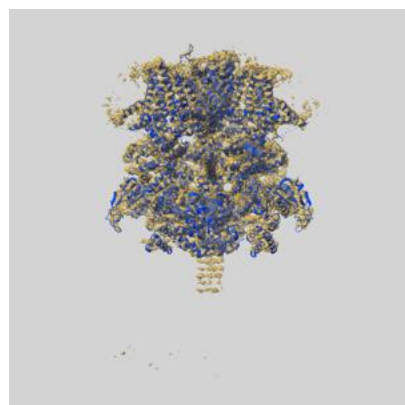
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.46	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.90	8.71	7.38

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.90 differs from the reported value 4.46 by more than 10 %

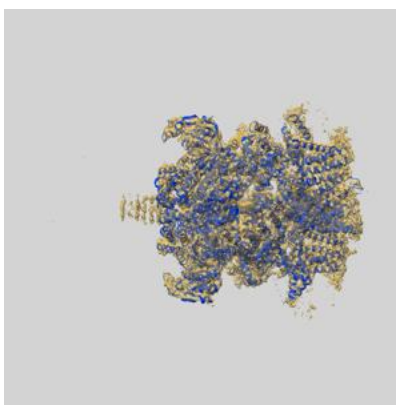
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-70215 and PDB model 9O82. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

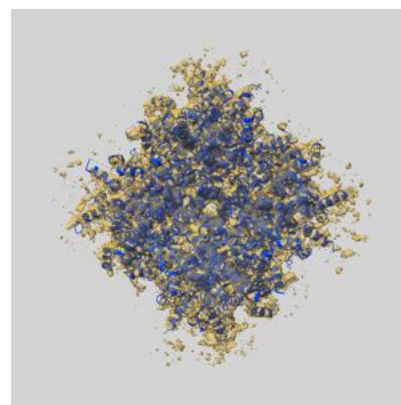
9.1 Map-model overlay [i](#)



X



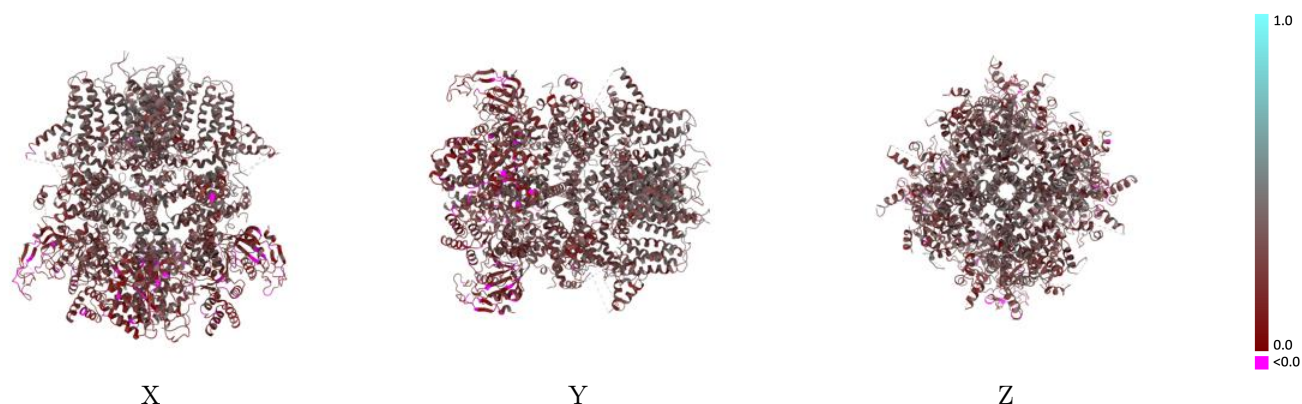
Y



Z

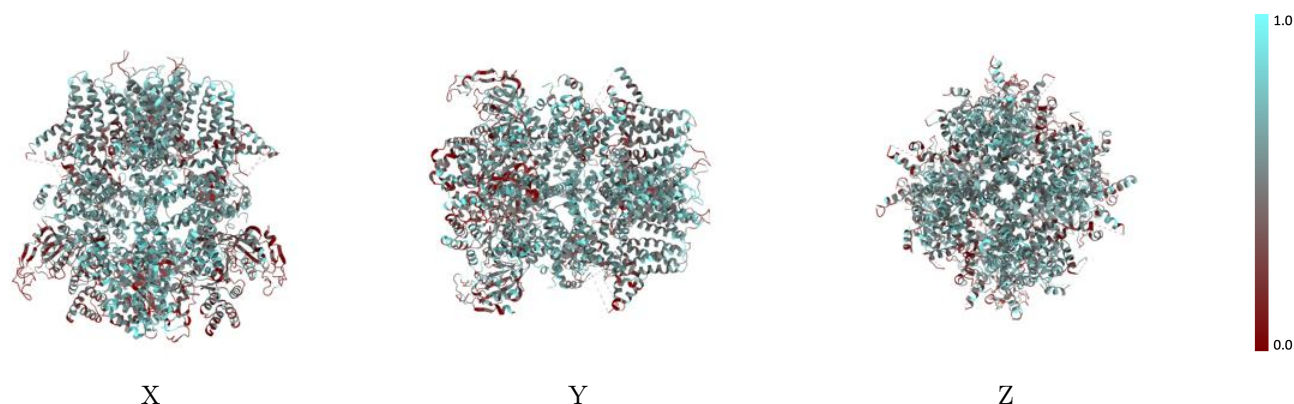
The images above show the 3D surface view of the map at the recommended contour level 0.0024 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



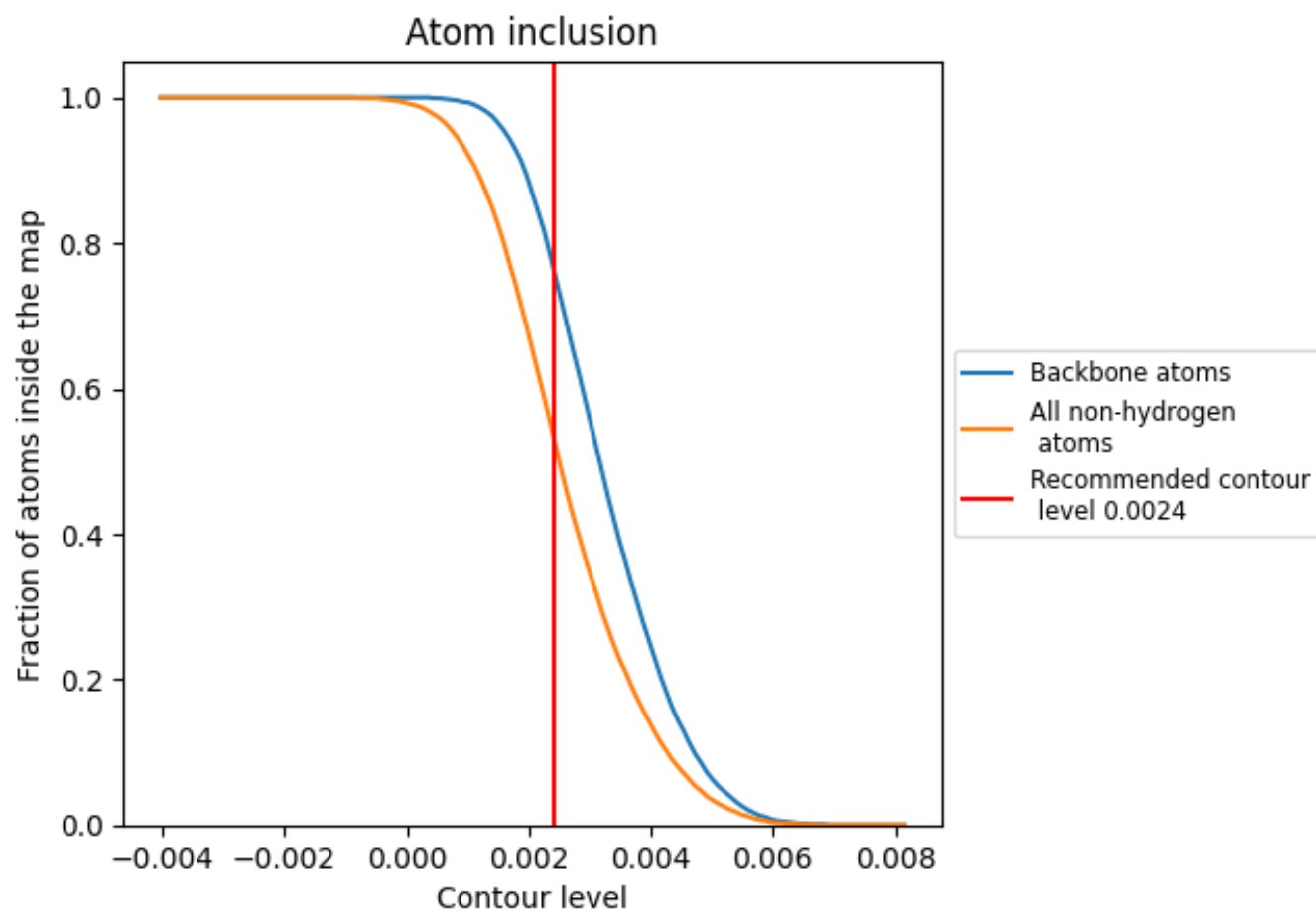
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0024).

9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 54% 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 (0.0024) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5350	0.2830
A	0.5530	0.2880
B	0.5130	0.2750
C	0.5560	0.2880
D	0.5170	0.2810

