



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 07:48 AM UTC

PDB ID : 8OR2 / pdb_00008or2
EMDB ID : EMD-17115
Title : CAND1-CUL1-RBX1-DCNL1
Authors : Shaaban, M.; Clapperton, J.A.; Ding, S.; Maeots, M.E.; Enchev, R.I.
Deposited on : 2023-04-12
Resolution : 3.20 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:

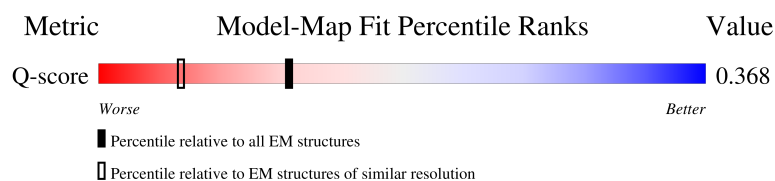
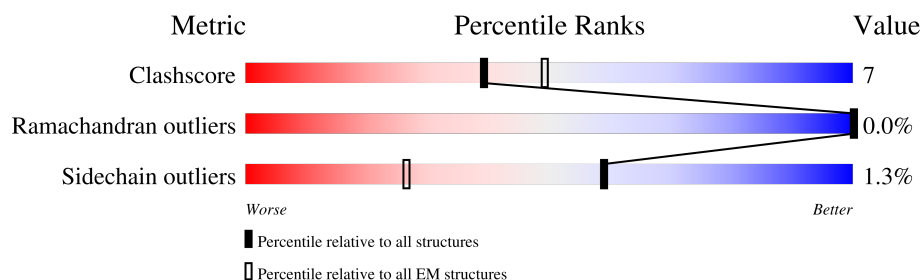
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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	15020 (2.70 - 3.70)

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	813	 6% 69% 13% 15%
2	B	108	 6% 66% 14% 19%
3	C	1238	 7% 82% 10% 9%
4	F	261	 7% 48% 10% 38%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16558 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 Cullin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	689	Total	C	N	O	S	0	0
			5651	3593	966	1064	28		

There are 37 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-36	MET	-	initiating methionine	UNP Q13616
A	-35	TRP	-	expression tag	UNP Q13616
A	-34	SER	-	expression tag	UNP Q13616
A	-33	HIS	-	expression tag	UNP Q13616
A	-32	PRO	-	expression tag	UNP Q13616
A	-31	GLN	-	expression tag	UNP Q13616
A	-30	PHE	-	expression tag	UNP Q13616
A	-29	GLU	-	expression tag	UNP Q13616
A	-28	LYS	-	expression tag	UNP Q13616
A	-27	GLY	-	expression tag	UNP Q13616
A	-26	SER	-	expression tag	UNP Q13616
A	-25	ALA	-	expression tag	UNP Q13616
A	-24	GLY	-	expression tag	UNP Q13616
A	-23	SER	-	expression tag	UNP Q13616
A	-22	ALA	-	expression tag	UNP Q13616
A	-21	ALA	-	expression tag	UNP Q13616
A	-20	GLY	-	expression tag	UNP Q13616
A	-19	SER	-	expression tag	UNP Q13616
A	-18	GLY	-	expression tag	UNP Q13616
A	-17	ALA	-	expression tag	UNP Q13616
A	-16	GLY	-	expression tag	UNP Q13616
A	-15	TRP	-	expression tag	UNP Q13616
A	-14	SER	-	expression tag	UNP Q13616
A	-13	HIS	-	expression tag	UNP Q13616
A	-12	PRO	-	expression tag	UNP Q13616
A	-11	GLN	-	expression tag	UNP Q13616
A	-10	PHE	-	expression tag	UNP Q13616
A	-9	GLU	-	expression tag	UNP Q13616

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	LYS	-	expression tag	UNP Q13616
A	-7	LEU	-	expression tag	UNP Q13616
A	-6	GLU	-	expression tag	UNP Q13616
A	-5	VAL	-	expression tag	UNP Q13616
A	-4	LEU	-	expression tag	UNP Q13616
A	-3	PHE	-	expression tag	UNP Q13616
A	-2	GLN	-	expression tag	UNP Q13616
A	-1	GLY	-	expression tag	UNP Q13616
A	0	PRO	-	expression tag	UNP Q13616

- Molecule 2 is a protein called E3 ubiquitin-protein ligase RBX1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	88	Total	C	N	O	S	0	0
			731	464	133	125	9		

- Molecule 3 is a protein called Cullin-associated NEDD8-dissociated protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	1132	Total	C	N	O	S	0	0
			8808	5608	1492	1652	56		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	GLY	-	expression tag	UNP Q86VP6
C	-6	SER	-	expression tag	UNP Q86VP6
C	-5	PRO	-	expression tag	UNP Q86VP6
C	-4	GLU	-	expression tag	UNP Q86VP6
C	-3	PHE	-	expression tag	UNP Q86VP6
C	-2	PRO	-	expression tag	UNP Q86VP6
C	-1	GLY	-	expression tag	UNP Q86VP6
C	0	ARG	-	expression tag	UNP Q86VP6

- Molecule 4 is a protein called DCN1-like protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	163	Total	C	N	O	S	0	0
			1365	889	225	245	6		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	GLY	-	expression tag	UNP Q96GG9
F	0	SER	-	expression tag	UNP Q96GG9

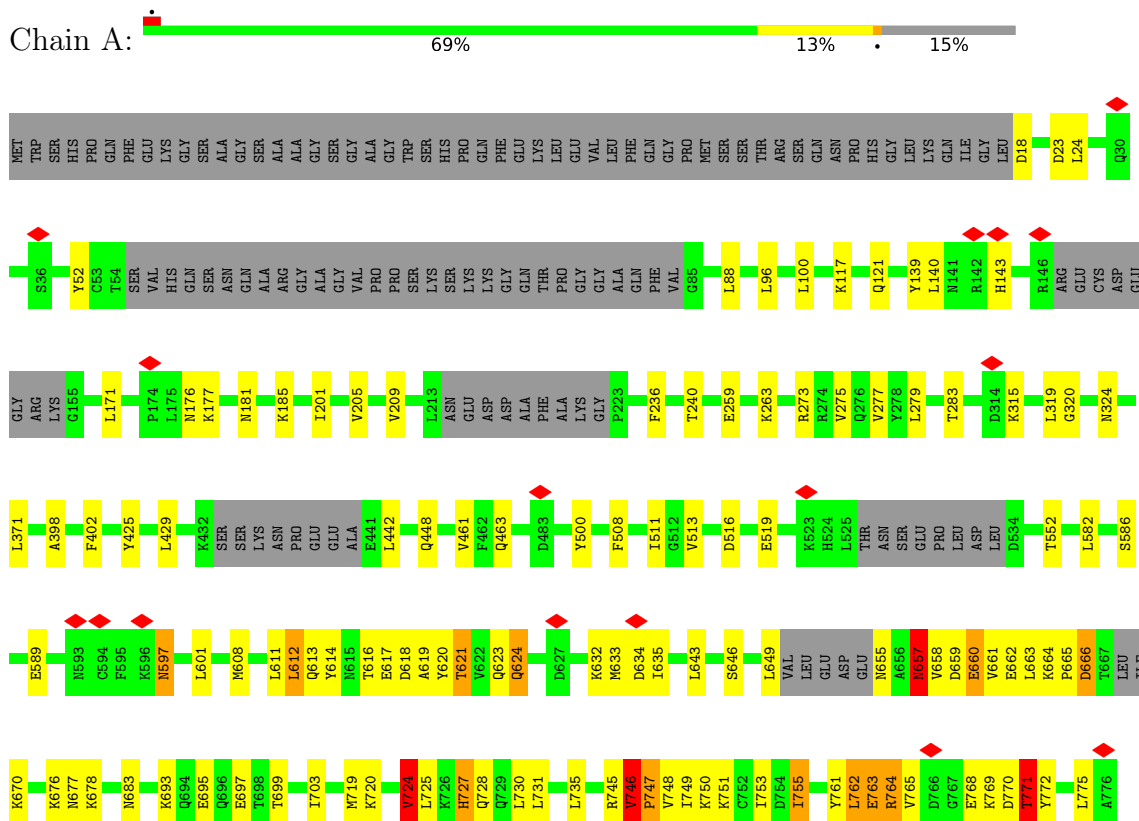
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
5	B	3	Total 3	Zn 3	0

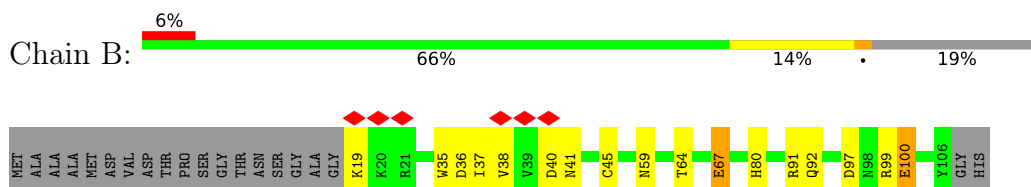
3 Residue-property plots

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.

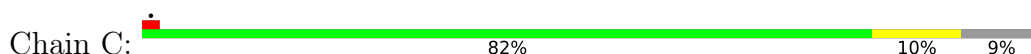
• Molecule 1: Cullin-1

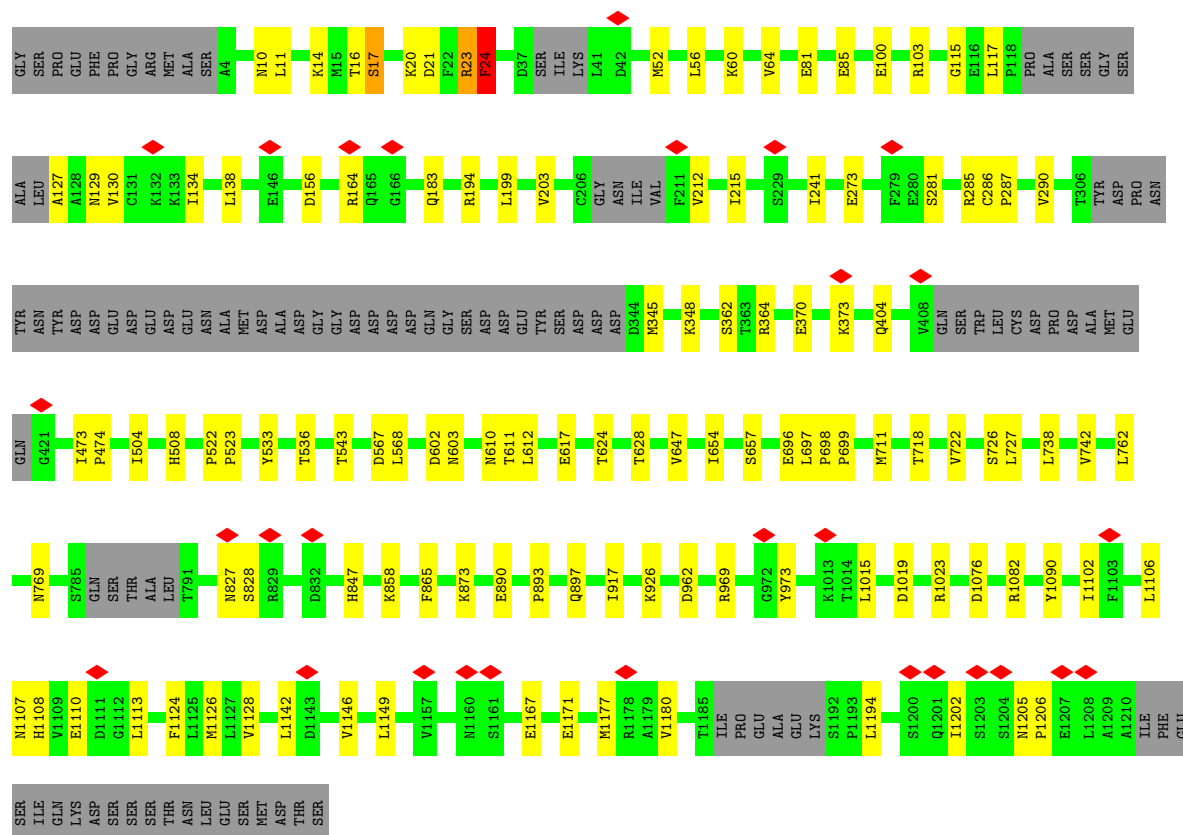


• Molecule 2: E3 ubiquitin-protein ligase RBX1

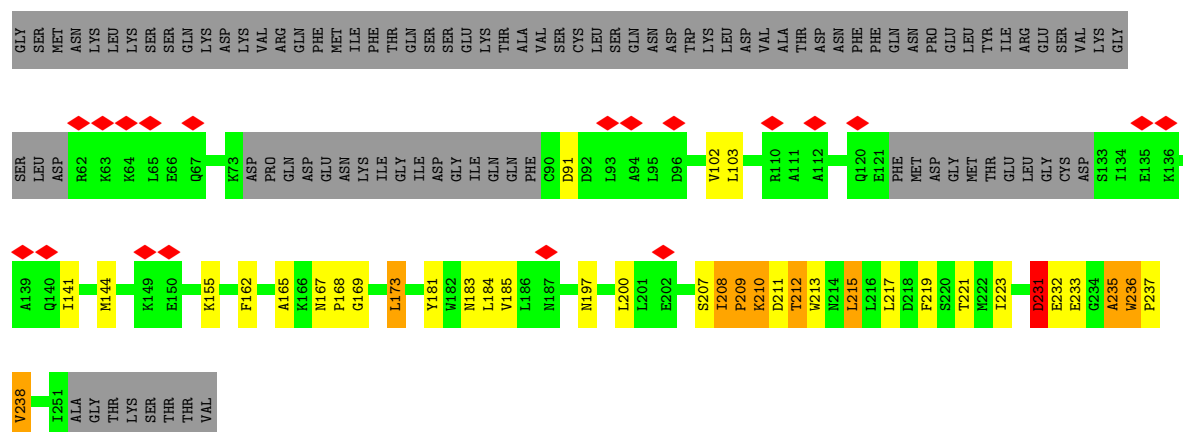


• Molecule 3: Cullin-associated NEDD8-dissociated protein 1





• Molecule 4: DCN1-like protein 1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	559316	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 ($e^-/\text{\AA}^2$)	49.1	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.037	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.0514	Depositor
Map size (Å)	345.6, 345.6, 345.6	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.59	13/5741 (0.2%)	0.54	15/7719 (0.2%)
2	B	0.42	2/752 (0.3%)	0.62	4/1020 (0.4%)
3	C	0.23	3/8943 (0.0%)	0.34	2/12107 (0.0%)
4	F	0.82	5/1399 (0.4%)	0.88	12/1883 (0.6%)
All	All	0.46	23/16835 (0.1%)	0.49	33/22729 (0.1%)

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	724	VAL	C-O	-7.69	1.15	1.24
1	A	755	ILE	CA-C	-7.60	1.43	1.52
2	B	91	ARG	C-N	7.54	1.45	1.33
1	A	720	LYS	C-O	-7.06	1.15	1.24
4	F	209	PRO	C-O	-6.98	1.15	1.23
1	A	764	ARG	CA-C	-6.82	1.44	1.52
1	A	755	ILE	CA-CB	-6.67	1.45	1.54
1	A	746	VAL	C-O	-6.56	1.16	1.24
4	F	233	GLU	N-CA	-6.32	1.42	1.46
3	C	21	ASP	C-O	-6.23	1.15	1.24
3	C	24	PHE	CA-CB	-6.19	1.45	1.54
3	C	23	ARG	C-O	-5.88	1.15	1.24
4	F	208	ILE	N-CA	-5.83	1.41	1.46
2	B	92	GLN	C-N	5.71	1.44	1.33
1	A	727	HIS	C-O	-5.69	1.17	1.24
1	A	755	ILE	C-O	-5.61	1.17	1.24
1	A	772	TYR	CA-CB	-5.56	1.41	1.53
4	F	209	PRO	CA-C	-5.53	1.45	1.52
1	A	727	HIS	CA-C	-5.53	1.45	1.52
4	F	219	PHE	CA-CB	-5.24	1.45	1.53
1	A	748	VAL	C-O	-5.22	1.18	1.24
1	A	768	GLU	C-O	-5.03	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	727	HIS	N-CA	-5.01	1.40	1.46

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	235	ALA	N-CA-C	10.08	125.11	113.01
1	A	658	VAL	N-CA-C	9.99	120.81	110.72
4	F	238	VAL	N-CA-C	9.97	125.45	112.98
3	C	17	SER	N-CA-C	9.48	124.11	112.54
2	B	92	GLN	O-C-N	-9.39	109.59	122.46
3	C	24	PHE	N-CA-C	-7.67	104.41	114.31
1	A	657	ASN	N-CA-C	7.51	121.05	108.13
1	A	666	ASP	N-CA-C	-7.33	98.04	109.25
2	B	92	GLN	CA-C-N	7.13	132.40	121.71
2	B	92	GLN	C-N-CA	7.13	132.40	121.71
4	F	238	VAL	CA-C-N	7.03	132.44	120.71
4	F	238	VAL	C-N-CA	7.03	132.44	120.71
1	A	751	LYS	CB-CA-C	-6.64	99.77	110.79
4	F	236	TRP	CA-C-N	6.05	127.41	119.84
4	F	236	TRP	C-N-CA	6.05	127.41	119.84
4	F	231	ASP	N-CA-C	-6.00	100.96	109.96
4	F	210	LYS	N-CA-C	5.98	117.79	111.28
1	A	772	TYR	N-CA-CB	-5.73	100.86	110.83
1	A	620	TYR	N-CA-C	5.67	116.86	108.86
4	F	208	ILE	CA-C-N	5.41	125.39	120.03
4	F	208	ILE	C-N-CA	5.41	125.39	120.03
1	A	755	ILE	CB-CA-C	-5.36	104.83	112.22
2	B	91	ARG	O-C-N	5.28	129.53	123.29
1	A	748	VAL	N-CA-C	5.27	116.04	110.72
4	F	169	GLY	N-CA-C	5.27	120.14	113.24
4	F	208	ILE	N-CA-CB	-5.26	106.80	112.37
1	A	624	GLN	N-CA-C	-5.20	105.69	111.36
1	A	762	LEU	N-CA-C	5.17	117.67	109.50
1	A	747	PRO	N-CA-C	-5.10	106.56	113.40
1	A	657	ASN	CA-C-N	5.05	127.60	120.42
1	A	657	ASN	C-N-CA	5.05	127.60	120.42
1	A	771	THR	N-CA-C	-5.05	99.39	108.48
1	A	728	GLN	N-CA-C	5.02	116.83	111.36

There are no chirality outliers.

There are no planarity outliers.

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	5651	0	5691	108	0
2	B	731	0	690	10	0
3	C	8808	0	9151	87	0
4	F	1365	0	1346	45	0
5	B	3	0	0	0	0
All	All	16558	0	16878	240	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (240) 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:727:HIS:CB	1:A:770:ASP:OD1	1.81	1.29
4:F:162:PHE:CE2	4:F:210:LYS:HB2	1.77	1.19
4:F:215:LEU:HD12	4:F:237:PRO:HD2	1.21	1.15
1:A:769:LYS:HE2	4:F:236:TRP:CZ3	1.84	1.10
1:A:727:HIS:HB2	1:A:770:ASP:OD1	1.48	1.07
4:F:162:PHE:HE2	4:F:210:LYS:HB2	0.94	1.07
1:A:719:MET:HG2	1:A:762:LEU:HD21	1.36	1.02
4:F:215:LEU:HD12	4:F:237:PRO:CD	1.88	1.02
4:F:162:PHE:HE2	4:F:210:LYS:CB	1.72	1.02
4:F:162:PHE:CE2	4:F:210:LYS:CB	2.43	1.02
4:F:215:LEU:CD1	4:F:237:PRO:HD2	1.92	0.97
1:A:618:ASP:O	1:A:670:LYS:CB	2.14	0.96
1:A:727:HIS:CG	1:A:770:ASP:OD1	2.24	0.91
1:A:727:HIS:HB3	1:A:770:ASP:OD1	1.72	0.89
1:A:769:LYS:HE2	4:F:236:TRP:CH2	2.07	0.87
4:F:215:LEU:CD1	4:F:237:PRO:CD	2.51	0.87
1:A:775:LEU:HD23	1:A:775:LEU:H	1.41	0.84
3:C:893:PRO:O	3:C:897:GLN:NE2	2.12	0.82
1:A:719:MET:CE	1:A:730:LEU:HD13	2.12	0.78
1:A:619:ALA:HB2	1:A:670:LYS:CB	2.13	0.78
1:A:657:ASN:ND2	1:A:657:ASN:O	2.17	0.77
1:A:745:ARG:O	1:A:749:ILE:HG13	1.85	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:345:MET:SD	3:C:348:LYS:NZ	2.57	0.77
1:A:181:ASN:OD1	1:A:185:LYS:NZ	2.18	0.76
1:A:761:TYR:O	1:A:775:LEU:HD23	1.86	0.75
1:A:236:PHE:O	1:A:240:THR:HG23	1.87	0.74
4:F:165:ALA:HB3	4:F:173:LEU:HD11	1.68	0.74
3:C:281:SER:OG	3:C:285:ARG:NH1	2.21	0.73
4:F:183:ASN:OD1	4:F:197:ASN:ND2	2.22	0.72
1:A:660:GLU:N	1:A:660:GLU:OE1	2.22	0.72
2:B:45:CYS:SG	2:B:80:HIS:CE1	2.83	0.71
3:C:1019:ASP:O	3:C:1023:ARG:NE	2.23	0.70
3:C:156:ASP:OD1	3:C:194:ARG:NH1	2.24	0.70
3:C:654:ILE:O	3:C:657:SER:OG	2.11	0.69
3:C:1076:ASP:OD2	3:C:1082:ARG:NH2	2.26	0.69
3:C:20:LYS:C	3:C:20:LYS:HD3	2.18	0.68
1:A:761:TYR:O	1:A:775:LEU:CD2	2.40	0.68
3:C:115:GLY:O	3:C:164:ARG:NH1	2.26	0.68
4:F:165:ALA:HB1	4:F:173:LEU:HD21	1.75	0.67
4:F:162:PHE:CE2	4:F:210:LYS:CG	2.78	0.67
1:A:552:THR:OG1	1:A:632:LYS:NZ	2.26	0.67
1:A:775:LEU:H	1:A:775:LEU:CD2	2.09	0.66
3:C:543:THR:HG23	3:C:568:LEU:HD22	1.76	0.66
3:C:610:ASN:OD1	3:C:611:THR:N	2.28	0.66
1:A:765:VAL:HG13	1:A:771:THR:HG22	1.77	0.65
3:C:100:GLU:OE1	3:C:103:ARG:NH2	2.29	0.65
3:C:567:ASP:OD1	3:C:568:LEU:N	2.30	0.65
3:C:858:LYS:NZ	3:C:890:GLU:OE2	2.19	0.65
3:C:718:THR:O	3:C:722:VAL:HG22	1.97	0.65
1:A:597:ASN:C	1:A:597:ASN:HD22	2.02	0.64
1:A:608:MET:O	1:A:612:LEU:HG	1.97	0.64
3:C:56:LEU:HD21	3:C:64:VAL:CG1	2.28	0.64
3:C:20:LYS:HD3	3:C:20:LYS:O	1.98	0.64
3:C:117:LEU:O	3:C:164:ARG:NH1	2.31	0.64
4:F:209:PRO:HG2	4:F:212:THR:HG23	1.80	0.64
3:C:624:THR:O	3:C:628:THR:OG1	2.14	0.63
2:B:100:GLU:OE1	3:C:60:LYS:HD2	1.98	0.63
1:A:746:VAL:N	1:A:747:PRO:CD	2.60	0.63
1:A:745:ARG:HB3	1:A:747:PRO:HD2	1.80	0.62
3:C:290:VAL:HG12	3:C:290:VAL:O	1.98	0.62
1:A:463:GLN:O	1:A:500:TYR:OH	2.14	0.62
1:A:764:ARG:HH21	4:F:235:ALA:HB1	1.63	0.62
3:C:56:LEU:HD21	3:C:64:VAL:HG12	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1194:LEU:HD12	3:C:1194:LEU:O	1.99	0.62
1:A:763:GLU:O	1:A:763:GLU:HG2	1.99	0.62
4:F:215:LEU:HD11	4:F:237:PRO:HD3	1.82	0.61
1:A:719:MET:HE1	1:A:730:LEU:HD13	1.82	0.61
4:F:162:PHE:CD2	4:F:210:LYS:HG3	2.35	0.61
1:A:621:THR:HG23	1:A:621:THR:O	2.00	0.61
1:A:719:MET:HG2	1:A:762:LEU:CD2	2.22	0.60
3:C:286:CYS:O	3:C:364:ARG:NH2	2.34	0.60
1:A:769:LYS:CE	4:F:236:TRP:CH2	2.83	0.60
1:A:731:LEU:O	1:A:735:LEU:HD23	2.02	0.60
1:A:616:THR:HG22	1:A:616:THR:O	2.02	0.60
4:F:215:LEU:CD1	4:F:237:PRO:HD3	2.31	0.60
1:A:18:ASP:N	1:A:52:TYR:HH	1.99	0.59
1:A:619:ALA:CB	1:A:670:LYS:CB	2.80	0.59
1:A:664:LYS:O	1:A:664:LYS:HG3	2.02	0.59
4:F:208:ILE:O	4:F:208:ILE:HD12	2.03	0.58
4:F:162:PHE:CE2	4:F:210:LYS:HG3	2.38	0.58
3:C:81:GLU:OE2	3:C:127:ALA:N	2.37	0.58
4:F:155:LYS:HD2	4:F:217:LEU:HD13	1.84	0.58
1:A:727:HIS:O	1:A:727:HIS:ND1	2.37	0.57
1:A:769:LYS:HE2	4:F:236:TRP:CE3	2.33	0.57
3:C:926:LYS:NZ	3:C:962:ASP:OD2	2.37	0.57
1:A:765:VAL:CG1	1:A:771:THR:HG22	2.35	0.57
1:A:695:GLU:OE2	1:A:699:THR:OG1	2.21	0.57
1:A:755:ILE:HG23	2:B:64:THR:HG22	1.87	0.56
3:C:183:GLN:N	3:C:183:GLN:OE1	2.38	0.56
1:A:614:TYR:CE2	1:A:649:LEU:CD1	2.88	0.56
4:F:162:PHE:CE2	4:F:210:LYS:CA	2.89	0.56
1:A:727:HIS:HB3	1:A:770:ASP:CG	2.30	0.55
4:F:162:PHE:CZ	4:F:210:LYS:HA	2.41	0.55
1:A:96:LEU:O	1:A:100:LEU:HD23	2.07	0.55
1:A:724:VAL:HG13	1:A:724:VAL:O	2.06	0.55
3:C:14:LYS:O	3:C:17:SER:HB3	2.07	0.55
3:C:827:ASN:OD1	3:C:828:SER:N	2.39	0.55
1:A:617:GLU:O	2:B:19:LYS:N	2.39	0.55
1:A:745:ARG:CB	1:A:747:PRO:HD2	2.36	0.55
1:A:775:LEU:HD23	1:A:775:LEU:O	2.07	0.55
4:F:155:LYS:HD2	4:F:217:LEU:CD1	2.38	0.54
1:A:597:ASN:C	1:A:597:ASN:ND2	2.66	0.54
1:A:619:ALA:CA	1:A:670:LYS:CB	2.86	0.54
4:F:232:GLU:OE2	4:F:232:GLU:HA	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:ASP:OD1	1:A:635:ILE:N	2.41	0.53
1:A:398:ALA:O	1:A:402:PHE:CD1	2.61	0.53
3:C:711:MET:HE2	3:C:711:MET:HA	1.90	0.52
3:C:56:LEU:HD23	3:C:56:LEU:O	2.09	0.52
1:A:693:LYS:O	1:A:697:GLU:OE1	2.27	0.52
1:A:619:ALA:HA	1:A:670:LYS:CB	2.39	0.52
1:A:513:VAL:O	1:A:516:ASP:OD1	2.29	0.51
2:B:35:TRP:HB2	2:B:38:VAL:HG12	1.93	0.51
1:A:88:LEU:HD11	1:A:140:LEU:HD21	1.93	0.51
3:C:1167:GLU:O	3:C:1171:GLU:HG3	2.10	0.51
3:C:612:LEU:HD21	3:C:647:VAL:HG21	1.94	0.50
1:A:176:ASN:OD1	1:A:177:LYS:N	2.45	0.50
1:A:643:LEU:O	1:A:646:SER:OG	2.23	0.50
1:A:201:ILE:O	1:A:205:VAL:HG23	2.12	0.50
1:A:724:VAL:O	1:A:724:VAL:HG22	2.10	0.50
3:C:602:ASP:OD1	3:C:603:ASN:N	2.44	0.50
2:B:97:ASP:OD2	2:B:99:ARG:NE	2.45	0.49
3:C:16:THR:O	3:C:16:THR:HG23	2.12	0.49
3:C:24:PHE:C	3:C:24:PHE:CD2	2.90	0.49
3:C:286:CYS:N	3:C:287:PRO:CD	2.76	0.49
4:F:167:ASN:HB2	4:F:168:PRO:CD	2.42	0.49
4:F:200:LEU:HD12	4:F:200:LEU:O	2.12	0.49
1:A:259:GLU:O	1:A:263:LYS:HG2	2.12	0.49
3:C:362:SER:O	3:C:404:GLN:NE2	2.45	0.49
3:C:1107:ASN:OD1	3:C:1108:HIS:N	2.44	0.49
1:A:727:HIS:CD2	1:A:770:ASP:OD1	2.65	0.49
3:C:1142:LEU:HD23	3:C:1142:LEU:C	2.38	0.49
1:A:139:TYR:O	1:A:143:HIS:ND1	2.41	0.48
4:F:223:ILE:HG23	4:F:223:ILE:O	2.11	0.48
3:C:273:GLU:OE1	3:C:273:GLU:N	2.45	0.48
2:B:67:GLU:O	2:B:67:GLU:HG3	2.03	0.48
3:C:24:PHE:HD2	3:C:24:PHE:O	1.96	0.48
1:A:727:HIS:ND1	1:A:727:HIS:C	2.72	0.48
4:F:184:LEU:HD23	4:F:184:LEU:C	2.38	0.48
3:C:1124:PHE:O	3:C:1128:VAL:HG23	2.14	0.48
1:A:320:GLY:O	1:A:324:ASN:ND2	2.46	0.48
2:B:36:ASP:OD1	2:B:37:ILE:N	2.47	0.48
1:A:371:LEU:HD22	1:A:448:GLN:HB3	1.96	0.48
1:A:621:THR:HG23	1:A:624:GLN:HB2	1.96	0.47
3:C:24:PHE:CD2	3:C:24:PHE:O	2.67	0.47
3:C:1090:TYR:HB2	3:C:1126:MET:HE3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:611:LEU:C	1:A:613:GLN:H	2.20	0.47
1:A:618:ASP:C	1:A:670:LYS:CB	2.86	0.47
3:C:199:LEU:O	3:C:203:VAL:HG23	2.14	0.47
3:C:696:GLU:N	3:C:696:GLU:OE1	2.48	0.47
4:F:162:PHE:CE2	4:F:210:LYS:HA	2.50	0.47
3:C:52:MET:HE2	3:C:52:MET:HA	1.96	0.47
3:C:203:VAL:HG21	3:C:241:ILE:HG13	1.96	0.47
4:F:141:ILE:HA	4:F:144:MET:HE2	1.97	0.47
3:C:20:LYS:C	3:C:20:LYS:CD	2.85	0.47
3:C:1202:ILE:O	3:C:1206:PRO:HD3	2.15	0.47
3:C:10:ASN:OD1	3:C:11:LEU:N	2.48	0.47
1:A:117:LYS:O	1:A:121:GLN:OE1	2.32	0.46
3:C:917:ILE:HD12	3:C:917:ILE:H	1.81	0.46
3:C:1142:LEU:O	3:C:1146:VAL:HG12	2.16	0.46
1:A:442:LEU:HD23	1:A:442:LEU:H	1.81	0.46
3:C:134:ILE:O	3:C:138:LEU:HD13	2.15	0.46
3:C:543:THR:CG2	3:C:568:LEU:HD22	2.45	0.46
4:F:213:TRP:CD1	4:F:213:TRP:C	2.92	0.46
1:A:727:HIS:CG	1:A:727:HIS:O	2.69	0.45
3:C:726:SER:O	3:C:727:LEU:HD22	2.16	0.45
1:A:725:LEU:HD23	1:A:730:LEU:HB2	1.96	0.45
1:A:23:ASP:O	1:A:24:LEU:HB3	2.17	0.45
4:F:91:ASP:N	4:F:91:ASP:OD2	2.49	0.45
4:F:208:ILE:H	4:F:208:ILE:HG13	1.52	0.45
3:C:1102:ILE:HG22	3:C:1106:LEU:HD13	1.98	0.45
3:C:926:LYS:HE3	3:C:926:LYS:HA	1.98	0.45
1:A:676:LYS:HA	1:A:676:LYS:HE3	1.99	0.45
1:A:761:TYR:C	1:A:775:LEU:HD21	2.41	0.45
3:C:24:PHE:C	3:C:24:PHE:HD2	2.25	0.45
3:C:847:HIS:C	3:C:847:HIS:HD1	2.25	0.45
4:F:207:SER:OG	4:F:208:ILE:N	2.50	0.45
1:A:623:GLN:OE1	1:A:623:GLN:HA	2.11	0.44
3:C:769:ASN:OD1	3:C:769:ASN:O	2.36	0.44
4:F:167:ASN:HB2	4:F:168:PRO:HD2	1.99	0.44
1:A:666:ASP:OD1	1:A:666:ASP:N	2.42	0.44
3:C:522:PRO:N	3:C:523:PRO:HD2	2.32	0.44
1:A:425:TYR:O	1:A:429:LEU:HD13	2.18	0.44
1:A:589:GLU:OE1	1:A:601:LEU:N	2.50	0.44
3:C:1110:GLU:HA	3:C:1113:LEU:HD23	1.98	0.44
1:A:755:ILE:O	1:A:755:ILE:HG22	2.18	0.44
3:C:129:ASN:OD1	3:C:130:VAL:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:533:TYR:O	3:C:536:THR:OG1	2.30	0.44
3:C:738:LEU:O	3:C:742:VAL:HG23	2.18	0.44
3:C:212:VAL:O	3:C:215:ILE:HG22	2.19	0.43
1:A:263:LYS:HD3	1:A:263:LYS:N	2.33	0.43
3:C:612:LEU:CD2	3:C:647:VAL:HG11	2.48	0.43
3:C:1149:LEU:HD23	3:C:1149:LEU:O	2.18	0.43
2:B:40:ASP:OD1	2:B:41:ASN:N	2.51	0.43
1:A:273:ARG:O	1:A:277:VAL:HG23	2.18	0.43
4:F:181:TYR:O	4:F:185:VAL:HG22	2.17	0.43
1:A:775:LEU:CD2	1:A:775:LEU:N	2.73	0.43
1:A:769:LYS:NZ	4:F:235:ALA:HB3	2.34	0.43
3:C:504:ILE:O	3:C:508:HIS:ND1	2.51	0.43
3:C:762:LEU:HD23	3:C:762:LEU:O	2.19	0.43
1:A:657:ASN:ND2	1:A:657:ASN:C	2.70	0.42
4:F:162:PHE:CD1	4:F:173:LEU:HD12	2.54	0.42
1:A:633:MET:HE2	1:A:633:MET:HA	2.01	0.42
3:C:865:PHE:O	3:C:873:LYS:NZ	2.46	0.42
1:A:719:MET:HE2	1:A:730:LEU:HD13	1.97	0.42
3:C:698:PRO:N	3:C:699:PRO:CD	2.81	0.42
3:C:1015:LEU:O	3:C:1015:LEU:HG	2.20	0.42
1:A:516:ASP:O	1:A:519:GLU:HG2	2.20	0.42
1:A:761:TYR:O	1:A:775:LEU:HD21	2.19	0.42
2:B:59:ASN:OD1	2:B:59:ASN:O	2.37	0.42
3:C:85:GLU:OE1	3:C:130:VAL:HG22	2.20	0.42
1:A:662:GLU:H	1:A:662:GLU:HG3	1.62	0.42
1:A:315:LYS:O	1:A:319:LEU:HD23	2.20	0.41
1:A:611:LEU:C	1:A:613:GLN:N	2.77	0.41
1:A:764:ARG:NH1	4:F:211:ASP:OD1	2.53	0.41
3:C:85:GLU:OE1	3:C:85:GLU:HA	2.20	0.41
3:C:969:ARG:O	3:C:973:TYR:CD2	2.73	0.41
1:A:275:VAL:HA	1:A:279:LEU:HB2	2.01	0.41
1:A:205:VAL:O	1:A:209:VAL:HG12	2.20	0.41
3:C:617:GLU:OE1	3:C:617:GLU:N	2.54	0.41
1:A:259:GLU:OE1	1:A:259:GLU:N	2.50	0.41
1:A:749:ILE:O	1:A:749:ILE:HG22	2.21	0.41
1:A:171:LEU:O	1:A:176:ASN:ND2	2.40	0.41
3:C:1177:MET:HE1	3:C:1205:ASN:HB3	2.03	0.41
1:A:461:VAL:HG22	1:A:703:ILE:HG21	2.02	0.41
1:A:683:ASN:OD1	1:A:683:ASN:C	2.64	0.41
3:C:473:ILE:N	3:C:474:PRO:CD	2.84	0.41
4:F:231:ASP:OD1	4:F:231:ASP:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:THR:O	1:A:283:THR:HG22	2.19	0.41
1:A:582:LEU:O	1:A:586:SER:OG	2.38	0.41
3:C:23:ARG:HH11	3:C:23:ARG:HD3	1.70	0.41
1:A:664:LYS:H	1:A:664:LYS:HG2	1.56	0.40
1:A:677:ASN:OD1	1:A:678:LYS:N	2.54	0.40
3:C:370:GLU:HA	3:C:373:LYS:HG2	2.04	0.40
3:C:1149:LEU:HD23	3:C:1149:LEU:C	2.46	0.40
1:A:508:PHE:O	1:A:511:ILE:HG22	2.21	0.40
3:C:697:LEU:N	3:C:698:PRO:CD	2.84	0.40
4:F:102:VAL:HG23	4:F:103:LEU:N	2.36	0.40
1:A:660:GLU:OE1	1:A:660:GLU:CA	2.70	0.40
3:C:1177:MET:HA	3:C:1180:VAL:HG22	2.03	0.40
3:C:1205:ASN:N	3:C:1206:PRO:CD	2.85	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	673/813 (83%)	654 (97%)	18 (3%)	1 (0%)	48	79
2	B	86/108 (80%)	84 (98%)	2 (2%)	0	100	100
3	C	1116/1238 (90%)	1076 (96%)	40 (4%)	0	100	100
4	F	157/261 (60%)	150 (96%)	7 (4%)	0	100	100
All	All	2032/2420 (84%)	1964 (97%)	67 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	612	LEU

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	625/725 (86%)	610 (98%)	15 (2%)	43	70
2	B	78/90 (87%)	76 (97%)	2 (3%)	40	69
3	C	1013/1104 (92%)	1012 (100%)	1 (0%)	88	91
4	F	146/234 (62%)	140 (96%)	6 (4%)	27	60
All	All	1862/2153 (86%)	1838 (99%)	24 (1%)	59	78

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	597	ASN
1	A	621	THR
1	A	655	ASN
1	A	657	ASN
1	A	659	ASP
1	A	660	GLU
1	A	661	VAL
1	A	663	LEU
1	A	665	PRO
1	A	724	VAL
1	A	746	VAL
1	A	750	LYS
1	A	753	ILE
1	A	763	GLU
1	A	771	THR
2	B	67	GLU
2	B	100	GLU
3	C	24	PHE
4	F	173	LEU
4	F	212	THR
4	F	215	LEU
4	F	221	THR
4	F	231	ASP
4	F	238	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	494	GLN
1	A	518	ASN
1	A	597	ASN
1	A	655	ASN
1	A	657	ASN
1	A	729	GLN
2	B	59	ASN
3	C	151	GLN
3	C	247	HIS
3	C	389	ASN
3	C	848	HIS
3	C	897	GLN
3	C	911	HIS
3	C	1035	HIS
4	F	159	GLN
4	F	170	GLN
4	F	214	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](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.

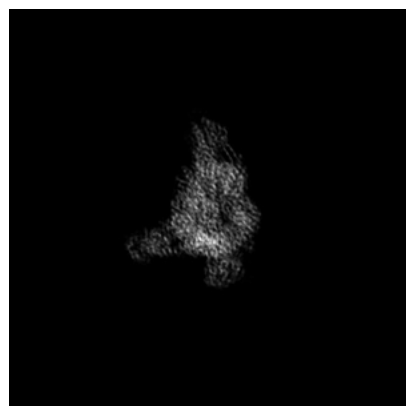
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17115. These allow visual inspection of the internal detail of the map and identification of artifacts.

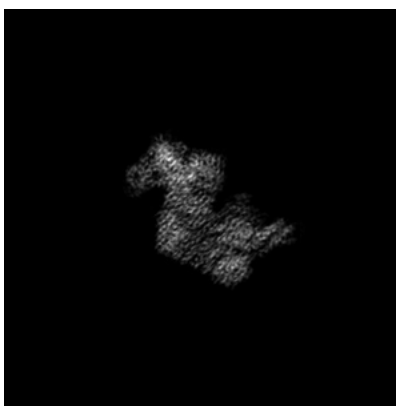
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

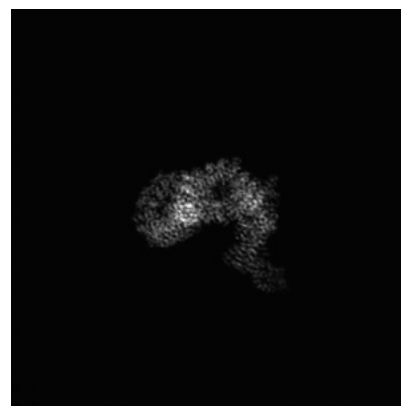
6.1.1 Primary map



X

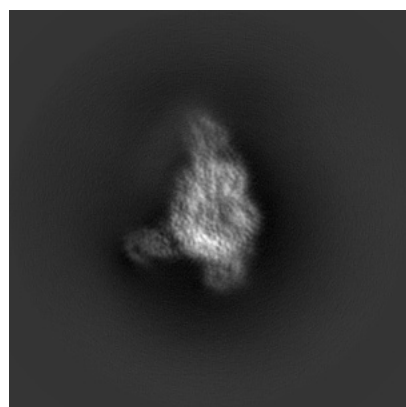


Y

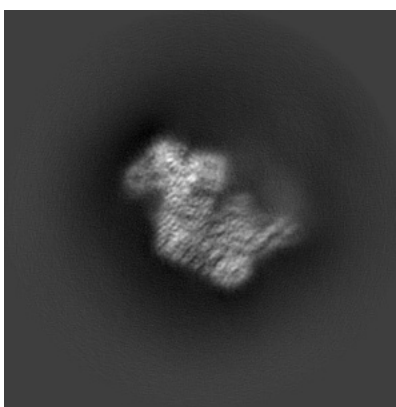


Z

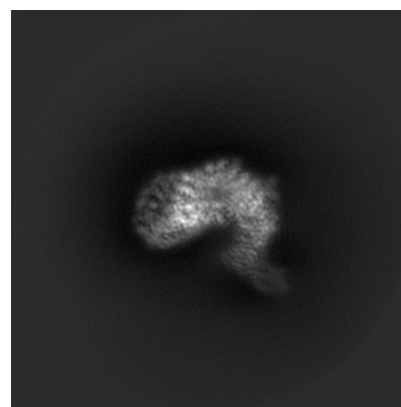
6.1.2 Raw map



X



Y

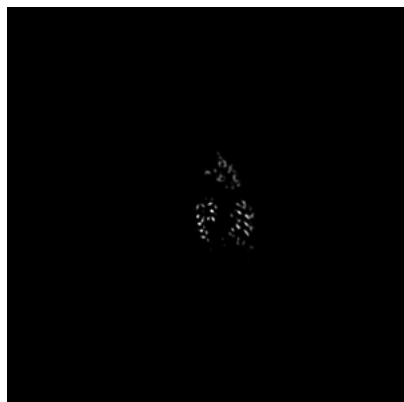


Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 160

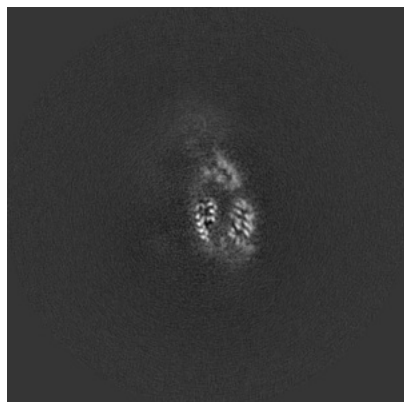


Y Index: 160

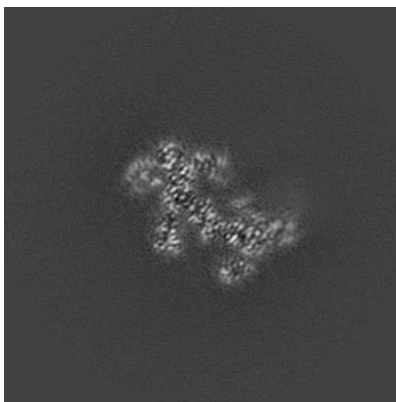


Z Index: 160

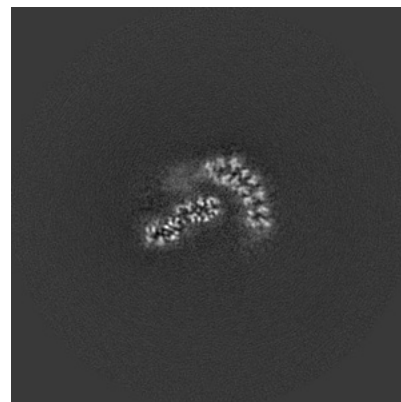
6.2.2 Raw map



X Index: 160



Y Index: 160



Z Index: 160

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: 141

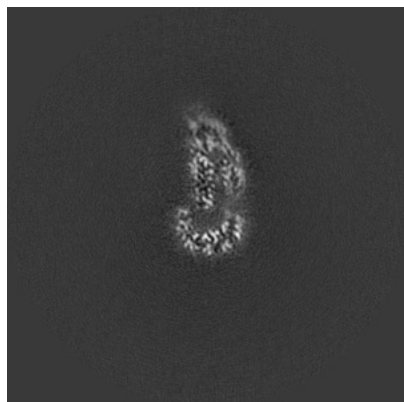


Y Index: 163

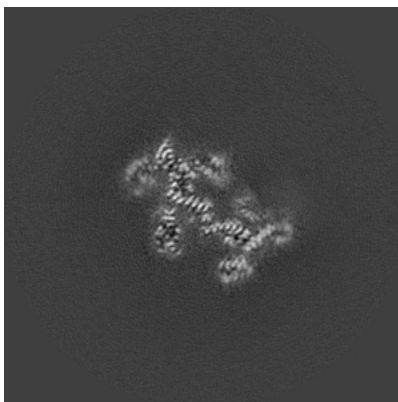


Z Index: 135

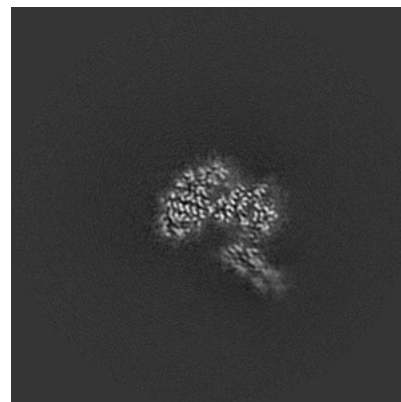
6.3.2 Raw map



X Index: 141



Y Index: 163

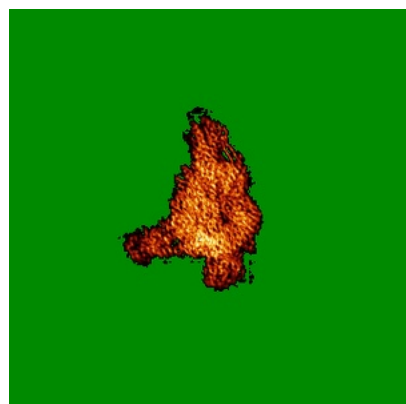


Z Index: 135

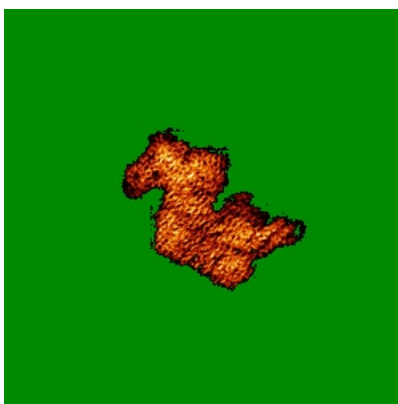
The images above show the largest variance slices of the map in three orthogonal directions.

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

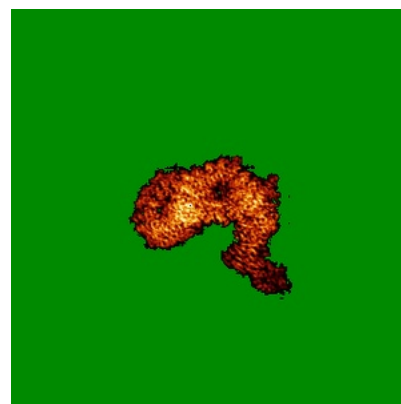
6.4.1 Primary map



X

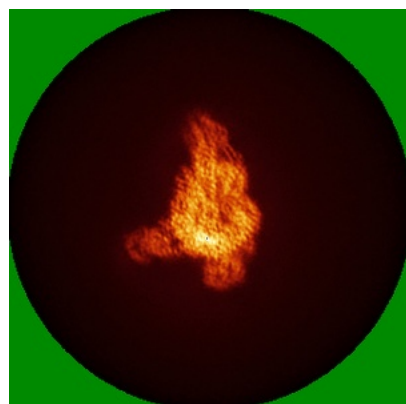


Y



Z

6.4.2 Raw map



X



Y

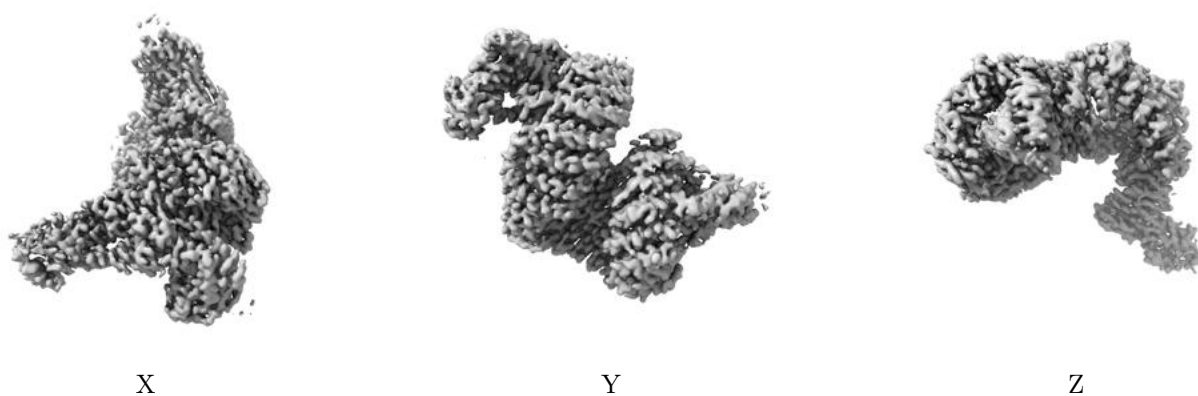


Z

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 0.0514. 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



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.

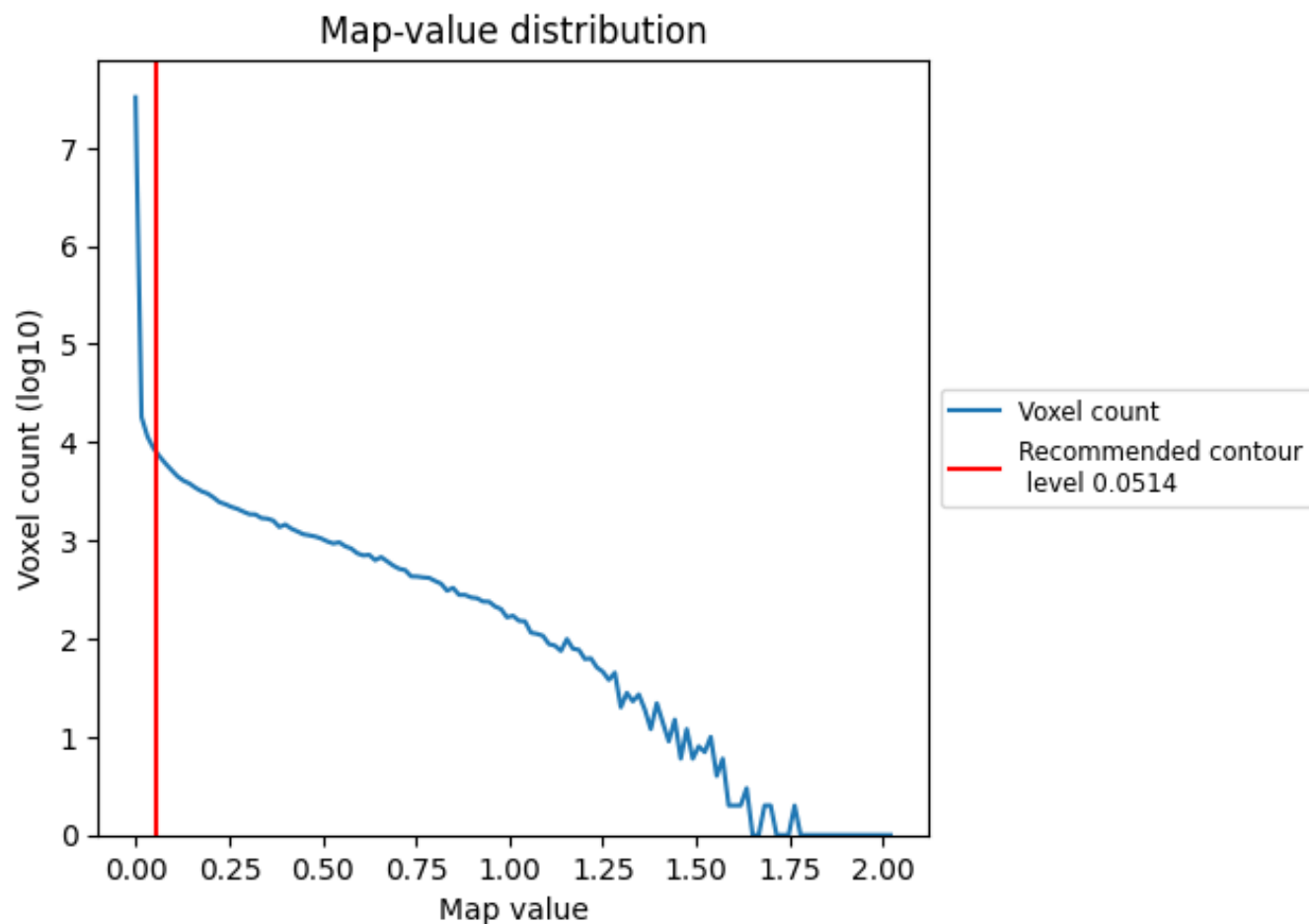
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

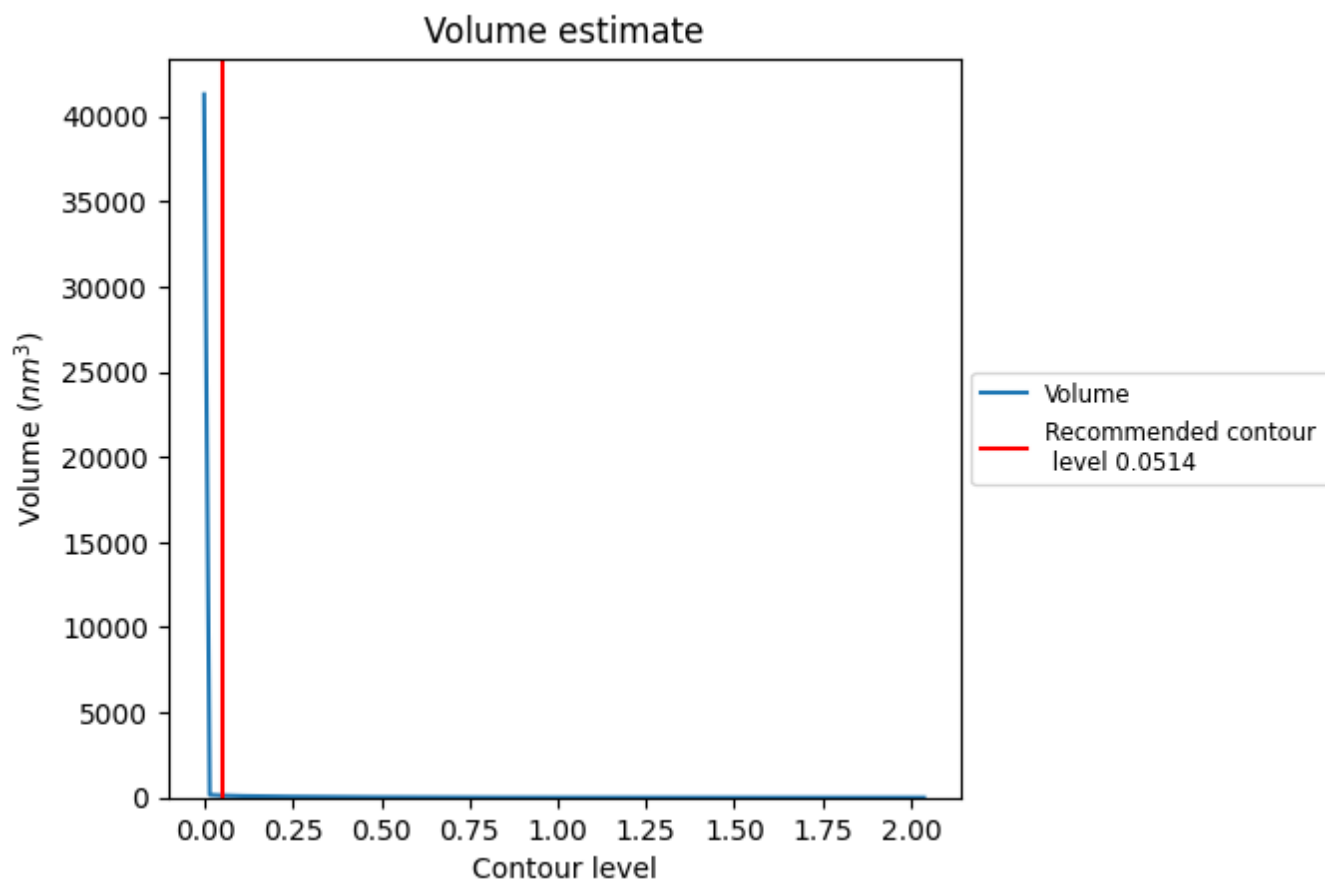
This section contains the results of statistical analysis of the map.

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.

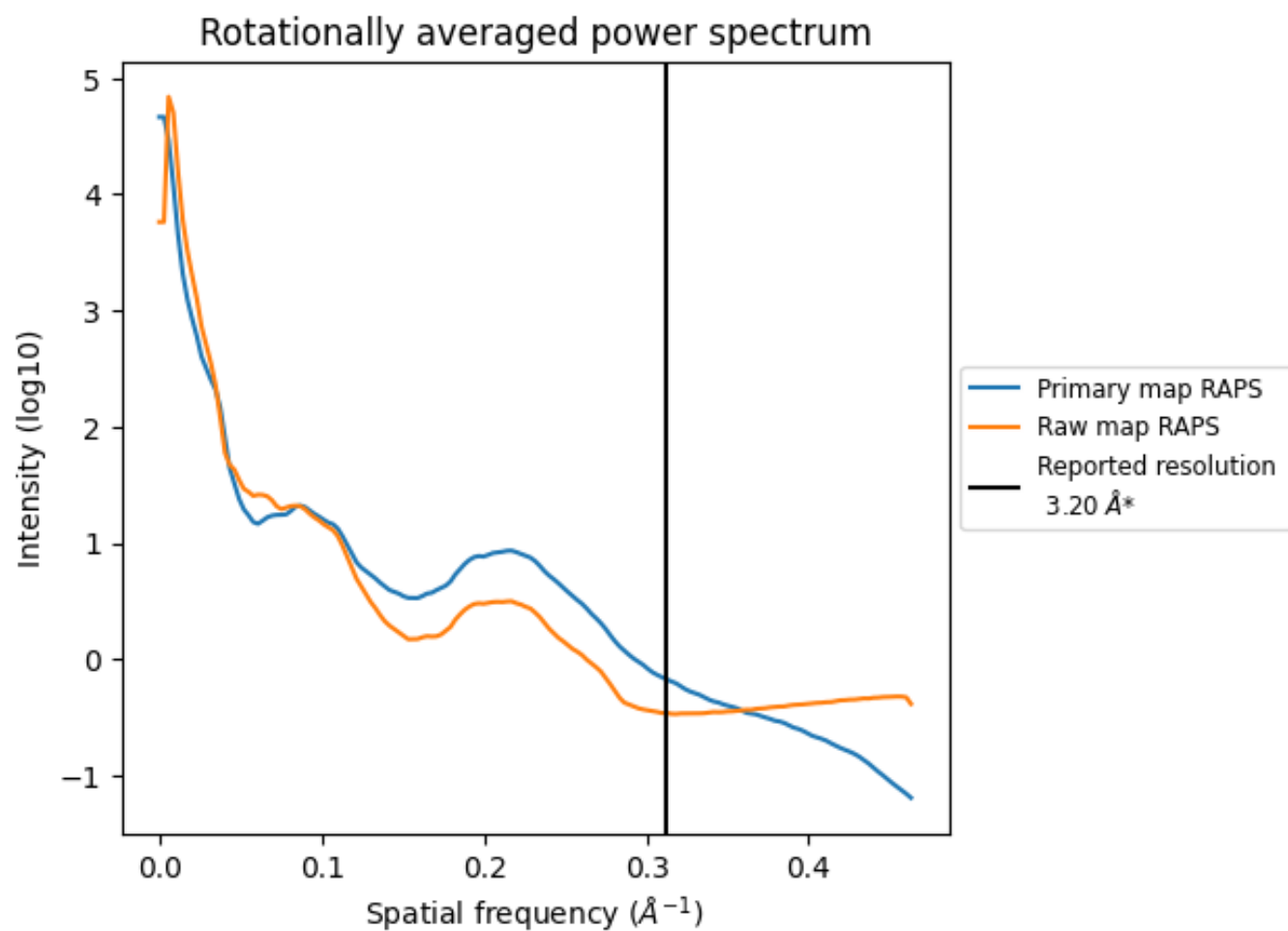
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 121 nm³; this corresponds to an approximate mass of 109 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 ⓘ

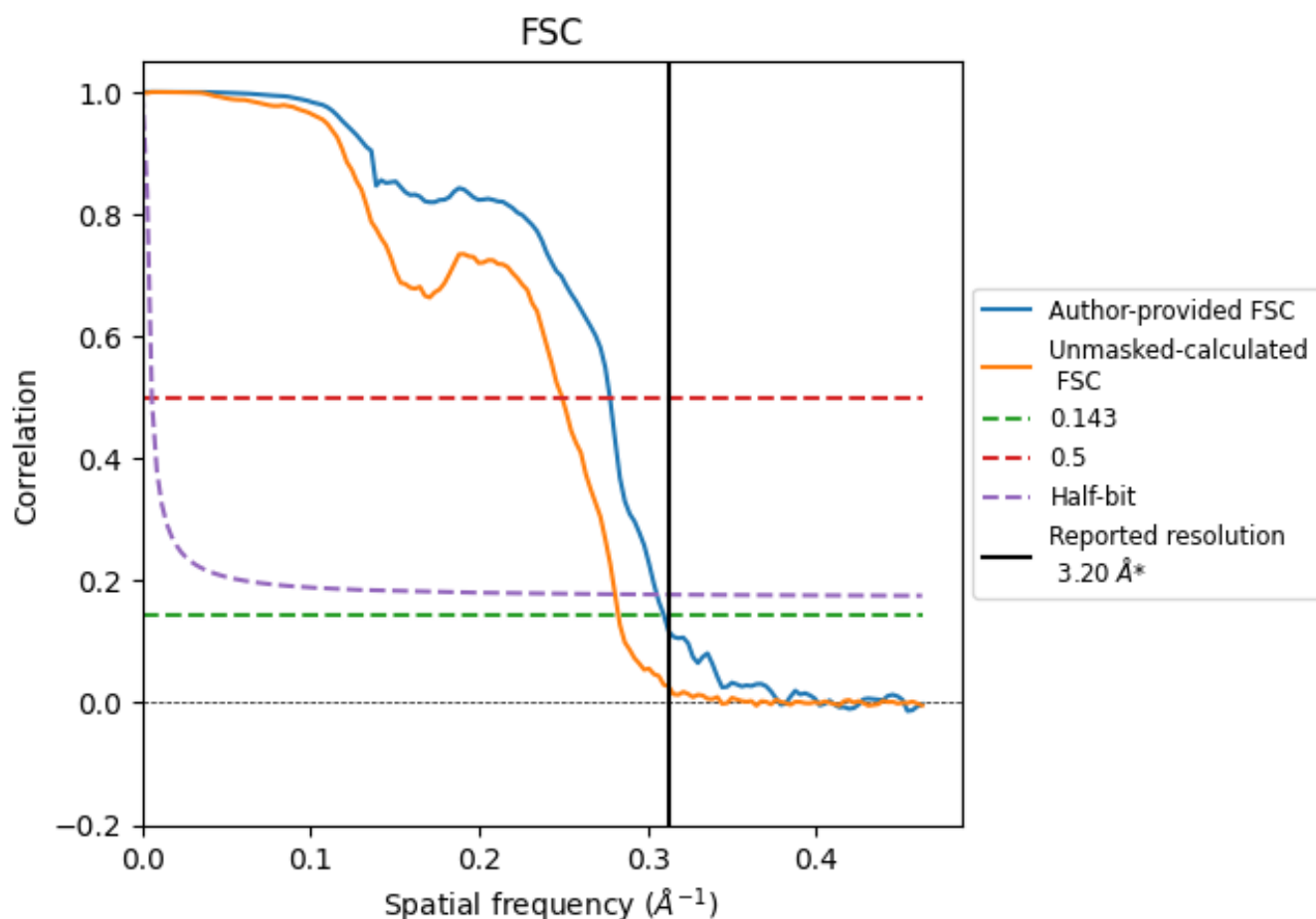


*Reported resolution corresponds to spatial frequency of 0.312 Å⁻¹

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 0.312 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

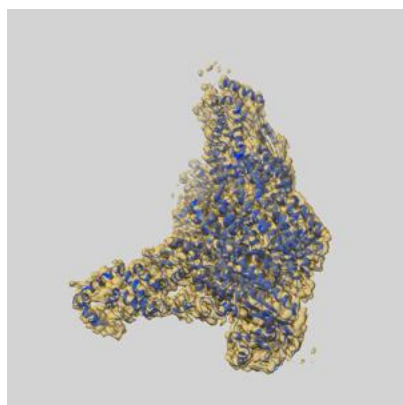
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.23	3.60	3.27
Unmasked-calculated*	3.54	4.01	3.56

*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 3.54 differs from the reported value 3.2 by more than 10 %

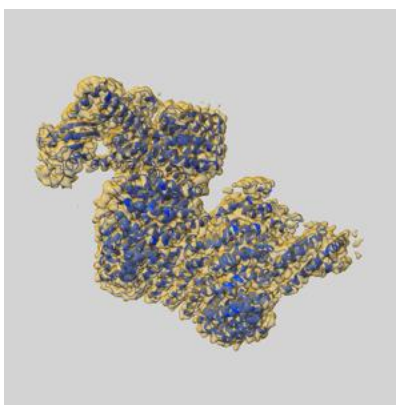
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-17115 and PDB model 8OR2. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

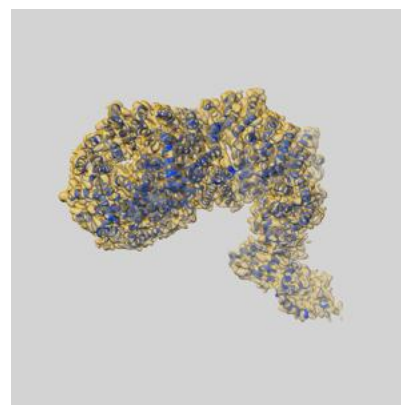
9.1 Map-model overlay [i](#)



X



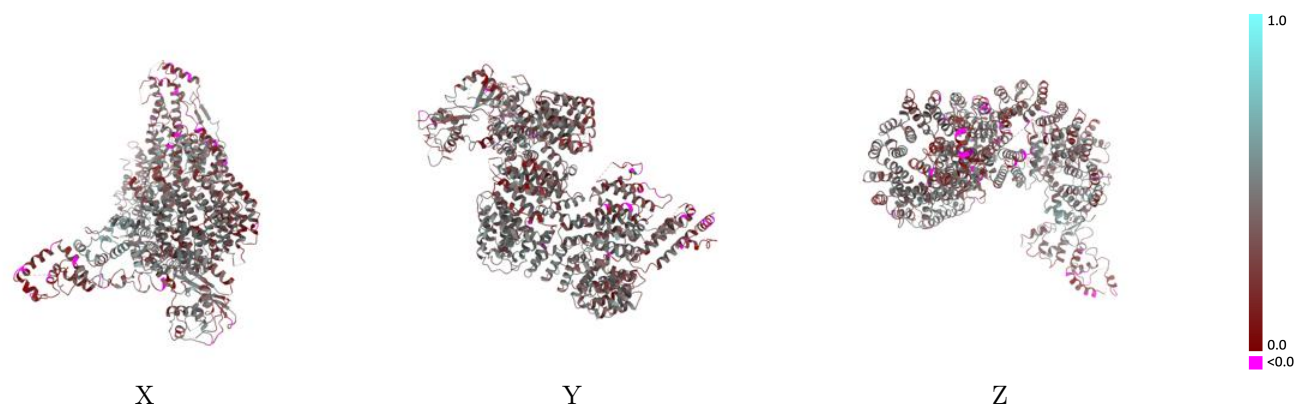
Y



Z

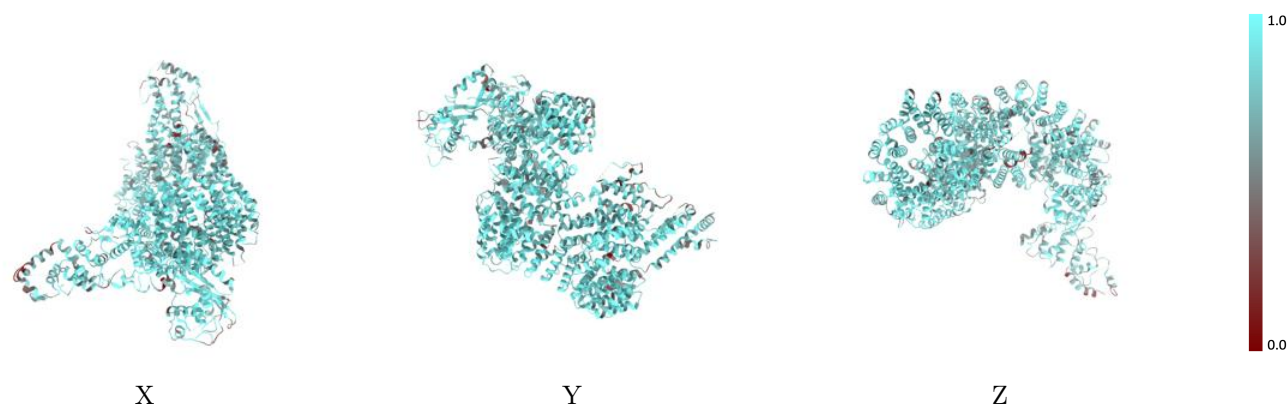
The images above show the 3D surface view of the map at the recommended contour level 0.0514 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



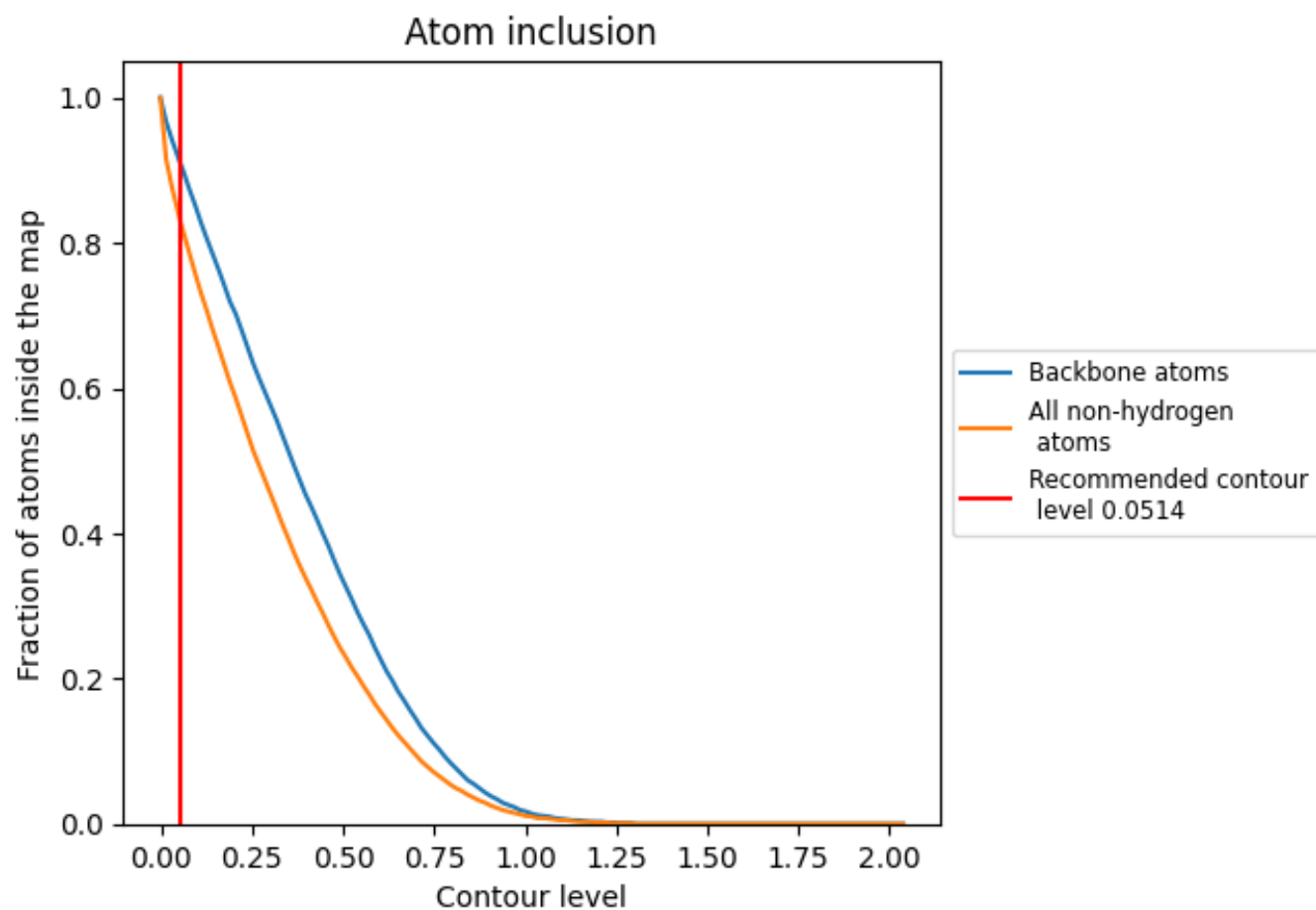
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0514).

9.4 Atom inclusion ⓘ



At the recommended contour level, 91% of all backbone atoms, 83% 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 (0.0514) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8330	0.3680
A	0.8420	0.3780
B	0.8200	0.3830
C	0.8480	0.3770
F	0.7050	0.2560

