



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 01:23 PM UTC

PDB ID : 5ZAK / pdb_00005zak
EMDB ID : EMD-6904
Title : Cryo-EM structure of human Dicer and its complexes with a pre-miRNA substrate
Authors : Liu, Z.; Wang, J.; Cheng, H.; Ke, X.; Sun, L.; Zhang, Q.C.; Wang, H.-W.
Deposited on : 2018-02-07
Resolution : 4.40 Å(reported)
Based on initial models : 5B16, 2KOU, 4GL2, 3C4T

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

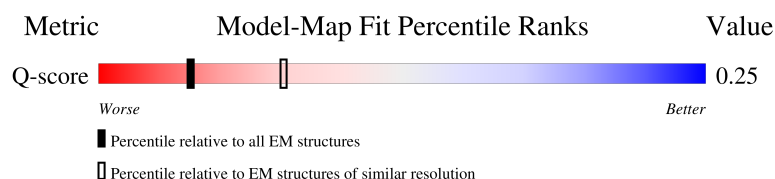
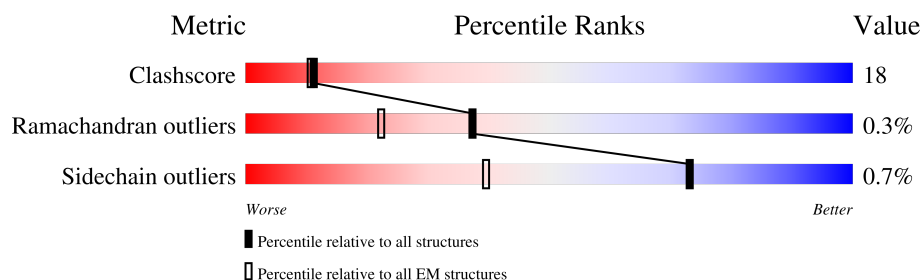
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

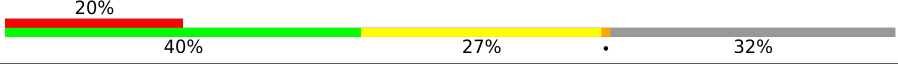
The reported resolution of this entry is 4.40 Å.

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	3132 (3.91 - 4.90)

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	1922	
2	B	366	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11112 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 Endoribonuclease Dicer.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1314	Total	C	N	O	S	0	0
			10552	6742	1807	1943	60		

- Molecule 2 is a protein called RISC-loading complex subunit TARBP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	75	Total	C	N	O	S	0	0
			560	345	99	110	6		



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	337566	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	2500	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	38462	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.116	Depositor
Minimum map value	-0.047	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.028	Depositor
Map size (Å)	219.49872, 219.49872, 219.49872	wwPDB
Map dimensions	168, 168, 168	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.30654, 1.30654, 1.30654	Depositor

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.42	0/10766	0.96	34/14558 (0.2%)
2	B	0.32	0/566	0.64	0/765
All	All	0.42	0/11332	0.95	34/15323 (0.2%)

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	13

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	134	THR	CA-C-N	9.03	137.95	121.70
1	A	134	THR	C-N-CA	9.03	137.95	121.70
1	A	116	HIS	CA-C-N	8.37	136.76	121.70
1	A	116	HIS	C-N-CA	8.37	136.76	121.70
1	A	649	LEU	CA-C-N	7.40	135.01	121.70
1	A	649	LEU	C-N-CA	7.40	135.01	121.70
1	A	137	ARG	CA-C-N	7.06	134.41	121.70
1	A	137	ARG	C-N-CA	7.06	134.41	121.70
1	A	1551	GLU	CA-C-N	7.01	134.32	121.70
1	A	1551	GLU	C-N-CA	7.01	134.32	121.70
1	A	862	LYS	N-CA-C	6.80	116.14	108.25
1	A	1591	LEU	N-CA-C	6.52	113.87	108.07
1	A	128	GLU	CA-C-N	5.95	132.41	121.70
1	A	128	GLU	C-N-CA	5.95	132.41	121.70
1	A	375	PRO	CA-C-N	5.87	132.75	121.54
1	A	375	PRO	C-N-CA	5.87	132.75	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	SER	CA-C-N	5.65	131.87	121.70
1	A	167	SER	C-N-CA	5.65	131.87	121.70
1	A	372	PHE	CA-C-N	5.62	131.82	121.70
1	A	372	PHE	C-N-CA	5.62	131.82	121.70
1	A	662	THR	N-CA-C	5.51	118.07	110.35
1	A	181	ILE	CA-C-N	5.49	129.06	120.82
1	A	181	ILE	C-N-CA	5.49	129.06	120.82
1	A	137	ARG	N-CA-C	5.45	120.83	114.62
1	A	366	ALA	N-CA-C	-5.38	107.37	114.04
1	A	452	TYR	N-CA-CB	-5.19	101.71	110.49
1	A	1591	LEU	N-CA-CB	-5.10	106.04	111.29
1	A	661	GLY	CA-C-N	5.09	128.90	121.31
1	A	661	GLY	C-N-CA	5.09	128.90	121.31
1	A	1588	LEU	N-CA-C	5.09	118.92	111.34
1	A	1610	PHE	CA-C-N	5.08	131.24	121.54
1	A	1610	PHE	C-N-CA	5.08	131.24	121.54
1	A	1308	ASP	N-CA-C	5.01	121.47	110.80
1	A	209	ILE	N-CA-C	-5.00	101.39	108.65

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1012	LEU	Peptide
1	A	131	ALA	Peptide
1	A	133	TRP	Peptide
1	A	1333	CYS	Peptide
1	A	1376	PRO	Peptide
1	A	140	GLN	Peptide
1	A	164	LEU	Peptide
1	A	165	SER	Peptide
1	A	1752	TYR	Peptide
1	A	207	ALA	Peptide
1	A	661	GLY	Peptide
1	A	707	GLU	Peptide
1	A	961	PRO	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	10552	0	10709	391	0
2	B	560	0	559	23	0
All	All	11112	0	11268	404	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 18.

All (404) 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:1882:VAL:O	1:A:1889:LYS:HA	1.65	0.96
1:A:769:TYR:O	1:A:805:LEU:HA	1.69	0.93
1:A:279:ASN:O	1:A:283:ASP:HB2	1.69	0.93
1:A:1313:GLU:O	1:A:1317:MET:HB2	1.74	0.88
1:A:1656:ASP:O	1:A:1659:LEU:HB2	1.80	0.82
1:A:543:TYR:HA	1:A:546:TYR:HB3	1.66	0.78
1:A:883:VAL:HA	1:A:889:THR:O	1.85	0.77
1:A:1908:ALA:O	1:A:1912:LEU:HB2	1.86	0.75
2:B:318:SER:HB3	2:B:322:LEU:H	1.52	0.74
1:A:1897:TYR:O	1:A:1901:LYS:HB2	1.87	0.74
1:A:543:TYR:HB3	1:A:579:TYR:HB3	1.68	0.74
1:A:1574:CYS:HB2	1:A:1578:ALA:HB2	1.70	0.73
1:A:443:CYS:SG	1:A:444:GLY:N	2.62	0.72
1:A:1771:ASP:O	1:A:1775:PHE:HB2	1.90	0.71
1:A:306:CYS:O	1:A:310:LEU:HB2	1.90	0.70
1:A:298:ILE:O	1:A:302:ILE:HB	1.92	0.69
1:A:654:ARG:H	1:A:667:LEU:HD23	1.59	0.67
1:A:336:GLU:HG3	1:A:338:GLU:H	1.60	0.67
1:A:271:LEU:HD13	1:A:310:LEU:HD23	1.76	0.66
1:A:384:LEU:HD11	1:A:561:ILE:HG21	1.77	0.66
1:A:1306:ALA:HB1	1:A:1372:SER:HB2	1.76	0.66
1:A:475:SER:HB2	1:A:497:GLU:H	1.60	0.66
1:A:1657:LYS:O	1:A:1661:HIS:N	2.28	0.66
1:A:372:PHE:H	1:A:373:VAL:HA	1.61	0.65
1:A:1353:ASN:O	1:A:1357:TYR:HB2	1.96	0.65
1:A:86:ARG:HA	1:A:94:LYS:HD3	1.77	0.65
1:A:449:GLU:HG2	1:A:450:ARG:HG3	1.79	0.65
1:A:67:GLY:HA3	1:A:70:LYS:HG2	1.79	0.64
1:A:640:ARG:O	1:A:672:ASN:ND2	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:LEU:HG	1:A:463:GLU:HB2	1.79	0.64
1:A:84:GLN:HG2	1:A:85:ILE:HD12	1.80	0.62
1:A:109:GLN:O	1:A:113:VAL:HB	1.98	0.62
1:A:382:GLU:O	1:A:386:LYS:N	2.31	0.62
1:A:470:GLU:HG3	1:A:516:LEU:HB3	1.82	0.62
1:A:635:ASN:HA	1:A:638:CYS:HB3	1.80	0.62
1:A:799:THR:HB	1:A:883:VAL:HB	1.82	0.62
1:A:1747:SER:O	1:A:1751:LYS:HB2	1.99	0.62
1:A:1757:TYR:O	1:A:1759:LYS:NZ	2.33	0.61
1:A:513:THR:HA	1:A:516:LEU:HB2	1.81	0.61
1:A:821:ARG:HH21	1:A:995:SER:HB2	1.64	0.61
1:A:257:CYS:HB3	1:A:567:ASP:HB3	1.83	0.61
1:A:1304:SER:HB2	1:A:1311:ASN:HD22	1.65	0.61
1:A:1839:ARG:HA	1:A:1842:ILE:HB	1.82	0.61
1:A:952:THR:HA	1:A:986:PRO:HB3	1.83	0.61
1:A:993:THR:HB	1:A:1048:TRP:HE1	1.66	0.61
1:A:651:PRO:HD2	1:A:670:PRO:HB3	1.80	0.61
1:A:633:HIS:O	1:A:636:ARG:HB3	2.01	0.61
1:A:71:THR:O	1:A:75:VAL:N	2.31	0.60
1:A:1588:LEU:HG	1:A:1589:LYS:HG3	1.81	0.60
1:A:920:PHE:HB3	1:A:982:ASN:HB2	1.82	0.60
1:A:114:ARG:HH21	1:A:123:GLU:H	1.49	0.60
1:A:1002:PRO:O	1:A:1006:ASN:N	2.27	0.60
1:A:831:LEU:HD11	1:A:834:SER:HB3	1.83	0.60
1:A:1004:HIS:O	1:A:1008:LYS:HB2	2.02	0.59
1:A:769:TYR:HB2	1:A:806:THR:O	2.02	0.59
1:A:640:ARG:HB2	1:A:670:PRO:HD2	1.84	0.59
1:A:279:ASN:O	1:A:283:ASP:CB	2.49	0.59
1:A:243:LEU:O	1:A:554:ARG:NH1	2.36	0.59
1:A:244:ASP:O	1:A:248:SER:N	2.35	0.59
1:A:51:LEU:HD22	1:A:62:VAL:HG22	1.84	0.59
1:A:134:THR:HB	1:A:135:LYS:HB2	1.84	0.59
1:A:1034:VAL:HG12	1:A:1036:GLU:H	1.68	0.58
1:A:1669:PHE:HZ	1:A:1757:TYR:H	1.50	0.58
1:A:473:TYR:HB3	1:A:501:GLN:HA	1.85	0.58
1:A:255:VAL:HG11	1:A:563:LEU:HD12	1.85	0.58
1:A:1026:SER:HA	1:A:1029:ASN:HB2	1.86	0.58
1:A:173:VAL:HG13	1:A:203:LEU:HB3	1.85	0.58
1:A:1621:CYS:SG	1:A:1694:TYR:N	2.72	0.58
1:A:1725:ARG:NH1	1:A:1846:SER:O	2.37	0.58
1:A:1332:PHE:HE1	1:A:1703:ARG:HB3	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1684:LEU:O	1:A:1688:THR:OG1	2.18	0.57
1:A:135:LYS:HA	1:A:138:TRP:HA	1.85	0.57
1:A:1583:LEU:O	1:A:1587:GLY:N	2.34	0.57
1:A:932:ILE:HB	1:A:1038:CYS:HB3	1.87	0.57
1:A:770:LEU:O	1:A:832:LYS:N	2.37	0.57
1:A:322:ALA:HA	1:A:352:LEU:HD22	1.86	0.57
1:A:1715:LEU:HD21	1:A:1827:LEU:HB3	1.87	0.57
1:A:1014:LEU:O	1:A:1019:LYS:NZ	2.38	0.56
1:A:246:TYR:OH	1:A:635:ASN:ND2	2.39	0.56
1:A:477:ASN:ND2	1:A:496:ALA:O	2.32	0.56
1:A:900:ILE:HG23	1:A:1042:PRO:HG2	1.87	0.56
2:B:325:CYS:SG	2:B:326:LEU:N	2.77	0.56
1:A:338:GLU:OE1	1:A:341:HIS:ND1	2.36	0.56
1:A:137:ARG:H	1:A:139:ASN:N	2.02	0.56
1:A:1692:TYR:HA	1:A:1761:VAL:HG23	1.88	0.56
1:A:863:PRO:HD2	1:A:869:PRO:HA	1.87	0.56
1:A:253:ILE:HD12	1:A:561:ILE:HG23	1.86	0.55
1:A:1732:LEU:HD23	1:A:1735:LEU:HD12	1.88	0.55
1:A:1834:TYR:HB3	1:A:1838:MET:HE2	1.88	0.55
1:A:510:ALA:O	1:A:514:ASN:ND2	2.40	0.55
1:A:86:ARG:NH2	1:A:87:GLY:O	2.40	0.55
1:A:365:PRO:HB3	1:A:368:LEU:HD13	1.87	0.55
1:A:380:LEU:O	1:A:384:LEU:HB2	2.07	0.55
1:A:1908:ALA:HA	1:A:1911:SER:HB3	1.89	0.55
1:A:283:ASP:OD2	2:B:324:GLN:NE2	2.40	0.54
1:A:702:LEU:O	1:A:706:GLY:N	2.40	0.54
1:A:1349:LYS:O	1:A:1355:ASN:ND2	2.34	0.54
1:A:1607:ARG:NH2	1:A:1682:TYR:OH	2.40	0.54
1:A:1614:GLN:NE2	1:A:1699:ASP:O	2.39	0.54
1:A:1019:LYS:O	1:A:1023:LYS:HB3	2.07	0.54
1:A:1359:LEU:HD11	1:A:1587:GLY:HA2	1.90	0.54
1:A:235:GLU:HA	1:A:590:CYS:HB3	1.89	0.54
1:A:771:TYR:HA	1:A:831:LEU:HA	1.90	0.54
1:A:1669:PHE:O	1:A:1672:LYS:HB3	2.08	0.54
1:A:175:ASP:HA	1:A:205:LEU:HB2	1.89	0.54
1:A:996:ARG:NH2	1:A:1001:THR:OG1	2.33	0.54
1:A:368:LEU:HD23	1:A:369:ASP:H	1.72	0.54
1:A:457:LEU:O	1:A:461:ILE:HB	2.07	0.54
1:A:800:ARG:HE	1:A:882:ASN:HB2	1.72	0.54
1:A:306:CYS:O	1:A:310:LEU:CB	2.56	0.54
1:A:1028:GLN:HA	1:A:1031:GLN:HE21	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1043:ILE:HG23	1:A:1047:LEU:HD23	1.90	0.54
1:A:92:ASN:H	1:A:95:ARG:HH22	1.56	0.53
1:A:770:LEU:HA	1:A:804:ILE:O	2.07	0.53
1:A:1708:GLY:HA2	1:A:1711:ILE:HD12	1.90	0.53
1:A:211:ASN:ND2	1:A:215:ASP:OD1	2.41	0.53
1:A:664:TYR:HA	1:A:680:VAL:HG11	1.90	0.53
1:A:1688:THR:HA	1:A:1759:LYS:HB2	1.90	0.53
1:A:354:LYS:HG3	2:B:361:ILE:HG21	1.89	0.53
1:A:931:VAL:HA	1:A:945:PHE:O	2.08	0.53
1:A:1353:ASN:O	1:A:1357:TYR:CB	2.57	0.53
1:A:1879:ARG:HD2	1:A:1891:LYS:HB3	1.91	0.53
1:A:372:PHE:N	1:A:373:VAL:HA	2.24	0.53
1:A:933:ILE:HD12	1:A:942:PRO:HG2	1.91	0.53
1:A:662:THR:HG22	1:A:685:SER:HB3	1.91	0.52
1:A:374:THR:HG21	1:A:378:ILE:H	1.74	0.52
1:A:680:VAL:HG12	1:A:682:PRO:HD3	1.92	0.52
1:A:70:LYS:O	1:A:73:ILE:HB	2.09	0.52
1:A:133:TRP:HB3	1:A:158:VAL:HG12	1.91	0.52
1:A:152:CYS:SG	1:A:153:TYR:N	2.83	0.52
1:A:190:MET:O	1:A:194:GLU:N	2.43	0.52
1:A:1688:THR:O	1:A:1695:ASN:ND2	2.42	0.52
1:A:1735:LEU:O	1:A:1739:LEU:HB2	2.10	0.52
1:A:132:SER:HB2	1:A:136:GLU:HG2	1.92	0.52
1:A:770:LEU:HD11	1:A:881:LEU:HD21	1.92	0.52
1:A:523:GLU:O	1:A:552:ARG:NH2	2.43	0.52
1:A:97:VAL:HA	1:A:147:VAL:HG13	1.91	0.52
1:A:375:PRO:HG3	1:A:440:ASN:HA	1.91	0.52
1:A:649:LEU:HB3	1:A:651:PRO:HD3	1.92	0.52
1:A:937:ARG:HH22	1:A:943:HIS:HB2	1.75	0.52
1:A:77:LEU:O	1:A:81:LEU:N	2.43	0.51
1:A:114:ARG:HG2	1:A:121:VAL:HG23	1.92	0.51
1:A:171:LEU:HD11	1:A:201:ARG:HB2	1.92	0.51
1:A:502:GLU:HA	1:A:505:LEU:HD12	1.92	0.51
1:A:1725:ARG:HH12	1:A:1847:ALA:HA	1.74	0.51
1:A:384:LEU:HD23	1:A:563:LEU:HD13	1.92	0.51
1:A:371:LYS:H	1:A:465:GLY:HA3	1.75	0.51
1:A:932:ILE:HG22	1:A:1040:ILE:HA	1.92	0.51
1:A:61:ILE:HB	1:A:234:ALA:HB2	1.93	0.51
1:A:769:TYR:HA	1:A:834:SER:HA	1.91	0.51
1:A:949:ASP:H	1:A:989:ASP:HB3	1.76	0.51
1:A:1070:LEU:HD13	1:A:1301:LEU:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1605:PRO:HG2	1:A:1825:MET:HE1	1.92	0.51
1:A:1656:ASP:HA	1:A:1659:LEU:HD13	1.93	0.50
1:A:299:SER:O	1:A:303:LEU:HB2	2.10	0.50
1:A:946:TYR:HD1	1:A:1045:ALA:HB1	1.76	0.50
1:A:1901:LYS:O	1:A:1905:ALA:CB	2.59	0.50
2:B:314:ILE:HD11	2:B:326:LEU:HG	1.92	0.50
1:A:113:VAL:HG12	1:A:121:VAL:HG21	1.92	0.50
1:A:225:LYS:HA	1:A:228:LYS:HE3	1.93	0.50
1:A:376:LYS:HG3	1:A:377:VAL:HG23	1.92	0.50
2:B:331:THR:OG1	2:B:333:PRO:O	2.21	0.50
1:A:259:PRO:HG3	1:A:385:ARG:HD2	1.94	0.50
1:A:343:LYS:HE2	2:B:312:LEU:HD22	1.93	0.50
1:A:1656:ASP:HB3	1:A:1659:LEU:HD22	1.94	0.50
1:A:134:THR:H	1:A:135:LYS:C	2.20	0.50
1:A:881:LEU:HA	1:A:891:ASP:O	2.12	0.50
1:A:1771:ASP:O	1:A:1775:PHE:CB	2.59	0.50
1:A:381:LEU:HA	1:A:384:LEU:HD13	1.94	0.49
1:A:1584:CYS:HA	1:A:1588:LEU:HB3	1.93	0.49
1:A:87:GLY:H	1:A:94:LYS:HB2	1.77	0.49
1:A:854:PHE:HA	1:A:858:LEU:HD12	1.95	0.49
1:A:466:LYS:HG2	1:A:468:ASP:H	1.77	0.49
1:A:246:TYR:HB3	1:A:554:ARG:HH22	1.77	0.49
2:B:329:LEU:N	2:B:335:THR:O	2.43	0.49
1:A:260:PHE:H	1:A:261:THR:HA	1.77	0.49
1:A:1009:GLY:HA2	1:A:1013:PRO:HG2	1.94	0.49
1:A:1697:ILE:HG22	1:A:1698:THR:HG23	1.95	0.49
1:A:90:SER:O	1:A:92:ASN:ND2	2.45	0.49
1:A:136:GLU:HB3	1:A:139:ASN:HA	1.95	0.49
1:A:451:ARG:HD3	1:A:484:ILE:HB	1.93	0.49
1:A:541:THR:OG1	1:A:542:GLU:N	2.46	0.49
2:B:350:GLU:OE1	2:B:353:ARG:NH2	2.39	0.49
1:A:447:PHE:N	1:A:535:VAL:O	2.46	0.49
1:A:442:LEU:HA	1:A:515:LEU:HD13	1.94	0.48
1:A:524:GLU:HA	1:A:552:ARG:HH22	1.78	0.48
1:A:73:ILE:O	1:A:76:LEU:HB2	2.13	0.48
1:A:194:GLU:HA	1:A:195:ASN:HA	1.63	0.48
1:A:92:ASN:O	1:A:95:ARG:NH1	2.46	0.48
1:A:128:GLU:HB3	1:A:130:ASN:H	1.76	0.48
1:A:1303:LEU:HD22	1:A:1369:MET:HE2	1.96	0.48
1:A:1660:ASN:HA	1:A:1663:ILE:HG12	1.96	0.48
2:B:289:ALA:HB3	2:B:357:GLN:HG3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:702:LEU:HD23	1:A:706:GLY:HA2	1.96	0.48
1:A:176:GLU:N	1:A:205:LEU:O	2.46	0.48
1:A:445:ILE:HG23	1:A:520:SER:H	1.78	0.48
1:A:457:LEU:HD11	1:A:519:THR:HG23	1.95	0.48
1:A:532:ASN:HB2	1:A:559:ASN:HD22	1.78	0.48
1:A:917:GLU:HG2	1:A:918:THR:HG23	1.96	0.48
1:A:1897:TYR:O	1:A:1901:LYS:CB	2.60	0.48
1:A:1807:ALA:HA	1:A:1810:ASP:HB2	1.96	0.48
1:A:667:LEU:N	1:A:678:SER:OG	2.47	0.48
1:A:1677:PHE:O	1:A:1680:LYS:NZ	2.45	0.48
1:A:65:ASN:OD1	1:A:544:ARG:NE	2.44	0.48
1:A:376:LYS:HA	1:A:380:LEU:HD22	1.96	0.48
1:A:454:ALA:HB1	1:A:471:LEU:HA	1.96	0.47
1:A:1908:ALA:O	1:A:1912:LEU:CB	2.59	0.47
1:A:312:VAL:HG12	1:A:313:LEU:HG	1.96	0.47
1:A:534:VAL:HG13	1:A:560:TYR:HA	1.96	0.47
1:A:252:GLU:HG3	1:A:560:TYR:HD2	1.79	0.47
1:A:449:GLU:OE2	1:A:450:ARG:NH1	2.48	0.47
1:A:181:ILE:O	1:A:187:ARG:NH2	2.42	0.47
1:A:1314:ARG:HB3	1:A:1721:TYR:HE1	1.80	0.47
1:A:214:CYS:SG	1:A:218:GLU:HG3	2.55	0.47
1:A:126:ASN:OD1	1:A:131:ALA:N	2.33	0.47
1:A:1340:GLU:O	1:A:1344:SER:CB	2.62	0.47
2:B:328:GLU:HG3	2:B:336:VAL:HG12	1.97	0.47
1:A:343:LYS:HD2	2:B:336:VAL:HG11	1.96	0.47
1:A:577:LYS:HA	1:A:580:LYS:HE2	1.96	0.47
1:A:671:ILE:HG22	1:A:674:PRO:HB3	1.96	0.47
1:A:680:VAL:HG13	1:A:694:VAL:HG11	1.96	0.47
1:A:1061:LEU:HD23	1:A:1064:LEU:HD12	1.96	0.47
1:A:1330:TYR:CZ	1:A:1588:LEU:HD22	2.49	0.47
2:B:313:ASP:OD2	2:B:348:ARG:NH2	2.48	0.47
1:A:320:LYS:HB2	1:A:456:VAL:HG22	1.96	0.47
1:A:1670:GLU:O	1:A:1674:ASN:N	2.45	0.47
1:A:862:LYS:HZ1	1:A:1374:PHE:HE1	1.61	0.47
1:A:508:PHE:HD1	1:A:513:THR:HG22	1.80	0.47
1:A:954:LEU:HB3	1:A:987:LEU:HD12	1.97	0.47
1:A:1864:THR:O	1:A:1885:VAL:N	2.44	0.47
1:A:82:SER:O	1:A:86:ARG:N	2.48	0.46
1:A:236:THR:OG1	1:A:590:CYS:SG	2.67	0.46
1:A:441:ILE:HG21	1:A:532:ASN:HD21	1.80	0.46
1:A:673:SER:HB2	1:A:708:LEU:HD22	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1717:THR:HG21	1:A:1736:ARG:HD2	1.97	0.46
1:A:536:ARG:HB3	1:A:539:LEU:HB3	1.95	0.46
1:A:1758:PHE:HD1	1:A:1811:ILE:HD13	1.81	0.46
1:A:53:GLU:HA	1:A:56:LEU:HD13	1.97	0.46
1:A:1365:LEU:HA	1:A:1368:ARG:HD3	1.96	0.46
1:A:1669:PHE:CZ	1:A:1756:LYS:HB2	2.51	0.46
1:A:258:GLY:O	1:A:260:PHE:HA	2.15	0.46
1:A:371:LYS:HA	1:A:371:LYS:HD3	1.70	0.46
1:A:643:SER:O	1:A:672:ASN:N	2.47	0.46
1:A:897:MET:HA	1:A:900:ILE:HG12	1.98	0.46
1:A:468:ASP:HB2	1:A:515:LEU:HG	1.98	0.46
1:A:496:ALA:HB1	1:A:501:GLN:HE22	1.79	0.46
1:A:571:SER:O	1:A:575:ASP:HB2	2.15	0.46
1:A:70:LYS:HD2	1:A:205:LEU:HB3	1.97	0.46
1:A:862:LYS:NZ	1:A:866:GLU:H	2.13	0.46
1:A:996:ARG:HE	1:A:999:LEU:HD23	1.81	0.46
1:A:1600:GLU:HA	1:A:1601:LYS:HA	1.62	0.46
1:A:118:ASP:OD2	1:A:1756:LYS:NZ	2.36	0.46
1:A:244:ASP:OD1	1:A:247:THR:OG1	2.33	0.46
1:A:1842:ILE:O	1:A:1846:SER:OG	2.30	0.46
1:A:137:ARG:HB3	1:A:140:GLN:H	1.80	0.46
1:A:923:LYS:HB3	1:A:925:GLU:HG2	1.96	0.46
1:A:128:GLU:HB3	1:A:130:ASN:N	2.31	0.46
1:A:1058:LEU:HD23	1:A:1061:LEU:HD12	1.98	0.46
1:A:137:ARG:HD3	1:A:140:GLN:HB2	1.99	0.45
1:A:257:CYS:HB2	1:A:568:LYS:HG3	1.98	0.45
1:A:779:THR:HB	1:A:824:GLU:HB3	1.98	0.45
1:A:271:LEU:HD11	1:A:307:ARG:HG3	1.97	0.45
1:A:446:ILE:H	1:A:519:THR:HG22	1.81	0.45
1:A:1601:LYS:HD3	1:A:1605:PRO:HG3	1.97	0.45
1:A:1755:HIS:HB3	1:A:1808:MET:HE1	1.98	0.45
1:A:1762:SER:HB3	1:A:1765:LEU:HD23	1.99	0.45
1:A:1823:SER:HB2	1:A:1826:SER:HB3	1.98	0.45
2:B:345:GLU:OE2	2:B:348:ARG:NH1	2.49	0.45
1:A:640:ARG:HG3	1:A:669:LEU:HB3	1.97	0.45
1:A:1852:SER:OG	1:A:1856:GLU:N	2.49	0.45
1:A:1340:GLU:O	1:A:1344:SER:HB3	2.16	0.45
1:A:1669:PHE:CE2	1:A:1757:TYR:HB2	2.52	0.45
1:A:1669:PHE:HE2	1:A:1757:TYR:HB2	1.82	0.45
1:A:1008:LYS:O	1:A:1012:LEU:N	2.42	0.45
1:A:1701:TYR:HE2	1:A:1810:ASP:HB3	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:308:HIS:N	2:B:330:SER:OG	2.47	0.45
1:A:121:VAL:HG12	1:A:147:VAL:HB	1.99	0.45
1:A:167:SER:HA	1:A:169:ILE:HG12	1.98	0.45
1:A:118:ASP:HB3	1:A:1668:ASN:HD21	1.82	0.45
1:A:128:GLU:H	1:A:129:VAL:HA	1.82	0.45
1:A:633:HIS:O	1:A:636:ARG:CB	2.65	0.45
1:A:1747:SER:O	1:A:1751:LYS:CB	2.63	0.45
1:A:295:SER:HB3	1:A:341:HIS:HD2	1.82	0.44
1:A:376:LYS:HB3	1:A:441:ILE:HG13	1.98	0.44
1:A:634:ILE:HD12	1:A:695:ALA:HB3	1.99	0.44
1:A:636:ARG:NH2	1:A:719:THR:OG1	2.50	0.44
1:A:900:ILE:HD12	1:A:1042:PRO:HG2	1.98	0.44
1:A:958:SER:O	1:A:968:PHE:N	2.50	0.44
1:A:1658:THR:HA	1:A:1661:HIS:HD2	1.81	0.44
1:A:347:PHE:CD1	2:B:337:CYS:HA	2.52	0.44
1:A:354:LYS:HD3	2:B:358:TYR:HD1	1.82	0.44
1:A:1315:LEU:HB2	1:A:1565:ALA:HB1	1.98	0.44
1:A:1612:SER:HB2	1:A:1615:LYS:HB2	1.99	0.44
1:A:256:ASP:HA	1:A:566:THR:HB	2.00	0.44
1:A:1612:SER:OG	1:A:1616:ASN:OD1	2.36	0.44
1:A:298:ILE:HG23	1:A:302:ILE:HD12	2.00	0.44
1:A:375:PRO:HB2	1:A:441:ILE:HB	1.99	0.44
1:A:1853:PRO:HB2	1:A:1905:ALA:HB3	1.98	0.44
1:A:319:ASP:HB2	1:A:359:CYS:HB2	1.98	0.44
1:A:229:ILE:O	1:A:231:LYS:NZ	2.49	0.44
1:A:931:VAL:HG21	1:A:1048:TRP:CG	2.53	0.44
1:A:111:SER:OG	1:A:112:ALA:N	2.51	0.44
1:A:853:ILE:HA	1:A:857:ILE:HD12	1.99	0.43
1:A:74:ALA:O	1:A:78:THR:N	2.52	0.43
1:A:359:CYS:HA	1:A:363:PHE:HD2	1.83	0.43
1:A:589:LYS:HA	1:A:592:LYS:HB2	1.99	0.43
1:A:956:PRO:HG2	1:A:983:LEU:HA	2.00	0.43
1:A:1557:LYS:HG3	1:A:1559:ILE:HG12	2.00	0.43
1:A:1855:ARG:HG2	1:A:1898:ARG:HH12	1.83	0.43
1:A:280:PHE:CE2	2:B:338:HIS:HB2	2.53	0.43
1:A:576:LEU:HA	1:A:579:TYR:HD2	1.83	0.43
1:A:932:ILE:HG13	1:A:945:PHE:HB2	2.01	0.43
1:A:369:ASP:HA	1:A:370:LEU:HA	1.81	0.43
1:A:240:LEU:HB2	1:A:244:ASP:HB2	2.00	0.43
1:A:462:LYS:N	1:A:469:PRO:HG3	2.33	0.43
1:A:53:GLU:HG2	1:A:654:ARG:HH22	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:486:LYS:HG3	1:A:521:ILE:HG13	1.99	0.43
1:A:899:ASP:HA	1:A:902:LYS:HD2	2.01	0.43
1:A:922:PHE:HE1	1:A:988:LEU:HD11	1.84	0.43
1:A:75:VAL:O	1:A:78:THR:OG1	2.20	0.43
1:A:447:PHE:CD1	1:A:523:GLU:HG2	2.54	0.43
1:A:651:PRO:HG3	1:A:668:TYR:HD2	1.84	0.43
1:A:667:LEU:HB2	1:A:678:SER:HB3	2.01	0.43
1:A:909:ILE:HD12	1:A:939:PHE:HA	2.01	0.43
1:A:286:ILE:HD13	2:B:314:ILE:HG23	2.01	0.42
1:A:536:ARG:NH2	1:A:539:LEU:O	2.52	0.42
1:A:1058:LEU:HA	1:A:1061:LEU:HB2	2.00	0.42
1:A:451:ARG:HD3	1:A:485:GLY:H	1.84	0.42
1:A:1688:THR:HG23	1:A:1759:LYS:HB2	2.00	0.42
1:A:1689:HIS:CD2	1:A:1691:SER:H	2.38	0.42
1:A:821:ARG:NH1	1:A:992:HIS:O	2.46	0.42
1:A:996:ARG:NH2	1:A:1006:ASN:OD1	2.52	0.42
1:A:1026:SER:O	1:A:1030:LYS:N	2.44	0.42
1:A:1339:HIS:H	1:A:1343:LEU:HD13	1.84	0.42
1:A:1594:ILE:H	1:A:1594:ILE:HG13	1.61	0.42
1:A:93:GLY:O	1:A:170:ASN:ND2	2.46	0.42
1:A:185:PRO:O	1:A:189:ILE:HG12	2.19	0.42
1:A:468:ASP:HB3	1:A:516:LEU:HA	2.01	0.42
1:A:153:TYR:HD1	1:A:156:LEU:HD21	1.84	0.42
1:A:989:ASP:HA	1:A:1032:ILE:HG23	2.00	0.42
1:A:992:HIS:HB3	1:A:1029:ASN:HD21	1.84	0.42
1:A:1047:LEU:HA	1:A:1050:LYS:HE2	2.02	0.42
1:A:333:ILE:HA	1:A:345:LEU:HD13	2.01	0.42
1:A:376:LYS:HD2	1:A:441:ILE:HG13	2.02	0.42
1:A:259:PRO:HA	1:A:260:PHE:CG	2.55	0.42
1:A:343:LYS:HZ2	2:B:328:GLU:HB2	1.85	0.42
1:A:240:LEU:O	1:A:244:ASP:CB	2.68	0.42
1:A:846:ILE:O	1:A:850:HIS:HB2	2.20	0.42
1:A:1819:ILE:O	1:A:1823:SER:N	2.51	0.42
1:A:225:LYS:HA	1:A:228:LYS:HG2	2.02	0.42
1:A:535:VAL:HG12	1:A:537:PHE:HB2	2.01	0.42
1:A:111:SER:O	1:A:1774:GLN:NE2	2.53	0.41
1:A:451:ARG:NH1	1:A:494:MET:O	2.41	0.41
1:A:925:GLU:HA	1:A:928:GLN:HG2	2.02	0.41
1:A:1873:THR:HG21	1:A:1879:ARG:HB3	2.01	0.41
1:A:631:ILE:HA	1:A:634:ILE:HB	2.02	0.41
1:A:1750:VAL:HG12	1:A:1773:VAL:HG22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:ARG:HG2	1:A:539:LEU:HD23	2.02	0.41
1:A:951:TYR:HD2	1:A:1032:ILE:HG21	1.84	0.41
1:A:373:VAL:HG23	1:A:466:LYS:HB2	2.02	0.41
1:A:631:ILE:O	1:A:635:ASN:N	2.40	0.41
1:A:1576:GLU:HG2	1:A:1580:GLN:HE21	1.85	0.41
1:A:1595:LYS:HA	1:A:1598:ASP:HB2	2.03	0.41
1:A:69:GLY:O	1:A:72:PHE:HB3	2.20	0.41
1:A:256:ASP:N	1:A:256:ASP:OD1	2.51	0.41
1:A:536:ARG:HD2	1:A:549:SER:HB2	2.03	0.41
1:A:1054:LEU:HD23	1:A:1057:ILE:HD12	2.02	0.41
1:A:1339:HIS:HB2	1:A:1342:ARG:HB3	2.02	0.41
1:A:77:LEU:HA	1:A:77:LEU:HD23	1.89	0.41
1:A:264:SER:HB3	1:A:267:TYR:HB3	2.02	0.41
1:A:364:SER:HA	1:A:365:PRO:HA	1.90	0.41
1:A:504:VAL:HA	1:A:507:LYS:HB2	2.03	0.41
1:A:792:LEU:HD22	1:A:822:SER:HA	2.02	0.41
1:A:865:LEU:HD22	1:A:1374:PHE:CD1	2.56	0.41
1:A:1316:GLU:O	1:A:1320:ASP:HB2	2.21	0.41
1:A:1817:GLY:O	1:A:1821:MET:HG2	2.21	0.41
1:A:64:LEU:HB3	1:A:67:GLY:HA2	2.03	0.41
1:A:73:ILE:HA	1:A:76:LEU:HD12	2.02	0.41
1:A:866:GLU:CD	1:A:1377:PRO:HB3	2.45	0.41
1:A:1863:GLU:HG2	1:A:1864:THR:HG23	2.02	0.41
1:A:1868:SER:O	1:A:1897:TYR:OH	2.24	0.41
1:A:1749:ALA:O	1:A:1754:TYR:N	2.46	0.41
1:A:283:ASP:HB3	2:B:324:GLN:HE22	1.85	0.41
1:A:893:ASP:HB3	1:A:897:MET:HE2	2.02	0.41
2:B:309:VAL:HA	2:B:329:LEU:HD23	2.03	0.41
1:A:65:ASN:ND2	1:A:544:ARG:HG2	2.36	0.41
1:A:316:TRP:CD1	1:A:320:LYS:HZ3	2.39	0.41
1:A:658:LEU:HD23	1:A:658:LEU:HA	1.91	0.41
1:A:1658:THR:O	1:A:1661:HIS:HB2	2.21	0.41
2:B:318:SER:OG	2:B:319:LEU:N	2.54	0.41
1:A:105:GLN:HA	1:A:108:GLN:HB2	2.03	0.40
1:A:274:LEU:HG	1:A:303:LEU:HD22	2.03	0.40
1:A:471:LEU:HD12	1:A:472:ALA:H	1.86	0.40
1:A:574:GLU:HA	1:A:577:LYS:HE2	2.01	0.40
1:A:781:LEU:HD12	1:A:782:PRO:HD2	2.03	0.40
1:A:1663:ILE:HA	1:A:1666:PHE:CD2	2.56	0.40
1:A:1839:ARG:O	1:A:1843:GLU:HB2	2.21	0.40
1:A:1758:PHE:CD1	1:A:1811:ILE:HD13	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:LYS:NZ	1:A:144:LYS:O	2.45	0.40
1:A:444:GLY:HA3	1:A:533:LEU:HB3	2.02	0.40
1:A:137:ARG:H	1:A:139:ASN:H	1.70	0.40
1:A:648:HIS:HA	1:A:649:LEU:HA	1.86	0.40
1:A:640:ARG:HH11	1:A:673:SER:HB3	1.86	0.40
1:A:1735:LEU:HD23	1:A:1735:LEU:HA	1.95	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	1296/1922 (67%)	1059 (82%)	233 (18%)	4 (0%)	36	71
2	B	73/366 (20%)	69 (94%)	4 (6%)	0	100	100
All	All	1369/2288 (60%)	1128 (82%)	237 (17%)	4 (0%)	37	71

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	875	ALA
1	A	139	ASN
1	A	659	PRO
1	A	259	PRO

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	1182/1730 (68%)	1175 (99%)	7 (1%)	78	80
2	B	60/301 (20%)	58 (97%)	2 (3%)	33	55
All	All	1242/2031 (61%)	1233 (99%)	9 (1%)	73	79

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	210	LEU
1	A	479	ILE
1	A	484	ILE
1	A	528	ILE
1	A	534	VAL
1	A	879	LEU
1	A	1882	VAL
2	B	336	VAL
2	B	360	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	146	GLN
1	A	493	GLN
1	A	532	ASN
1	A	559	ASN
1	A	635	ASN
1	A	1029	ASN
1	A	1299	GLN
1	A	1311	ASN
1	A	1325	HIS
1	A	1614	GLN
1	A	1660	ASN
1	A	1661	HIS
1	A	1668	ASN
1	A	1774	GLN
2	B	324	GLN

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-6904. These allow visual inspection of the internal detail of the map and identification of artifacts.

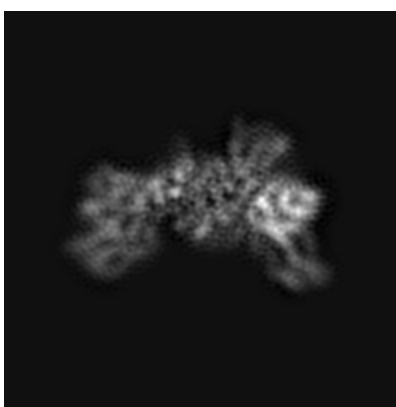
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

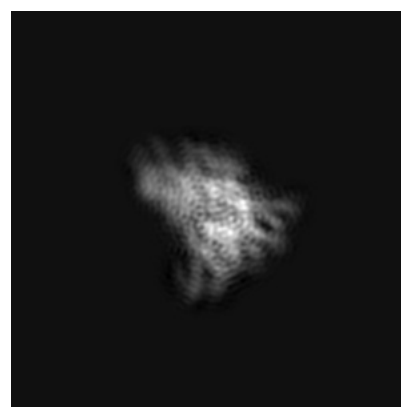
6.1.1 Primary map



X



Y

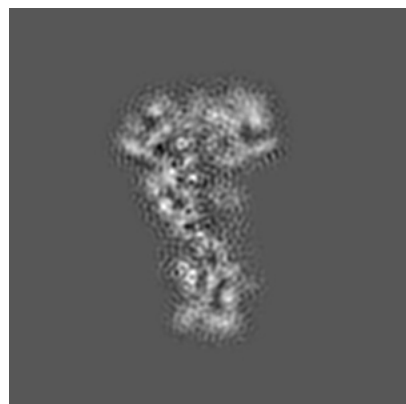


Z

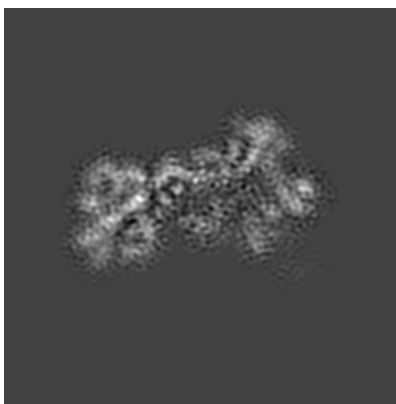
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 84



Y Index: 84



Z Index: 84

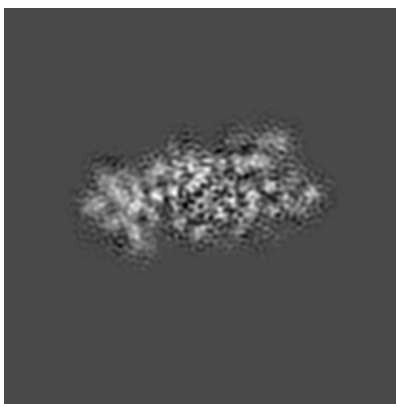
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

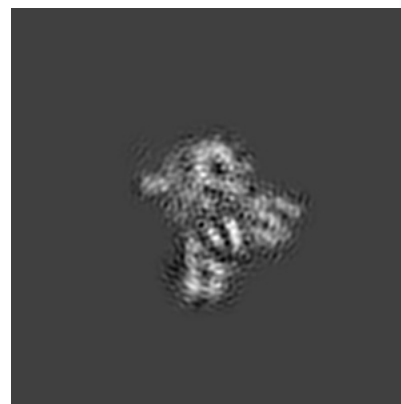
6.3.1 Primary map



X Index: 85



Y Index: 75



Z Index: 111

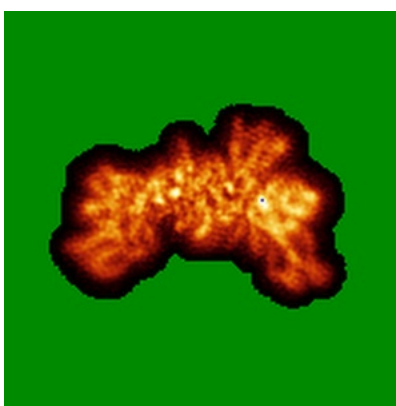
The images above show the largest variance slices of the map in three orthogonal directions.

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

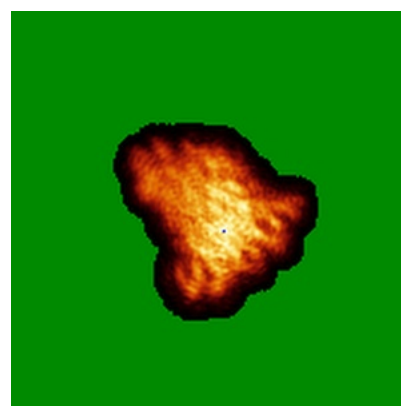
6.4.1 Primary map



X



Y

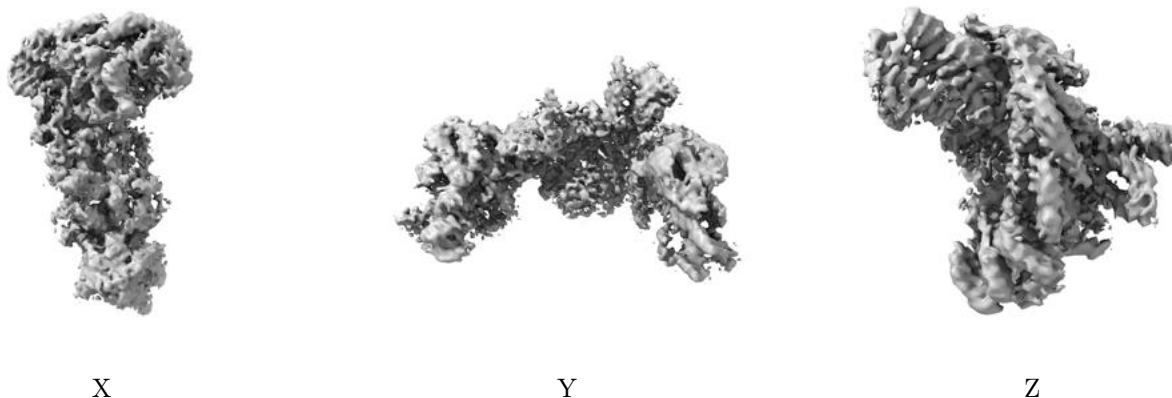


Z

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

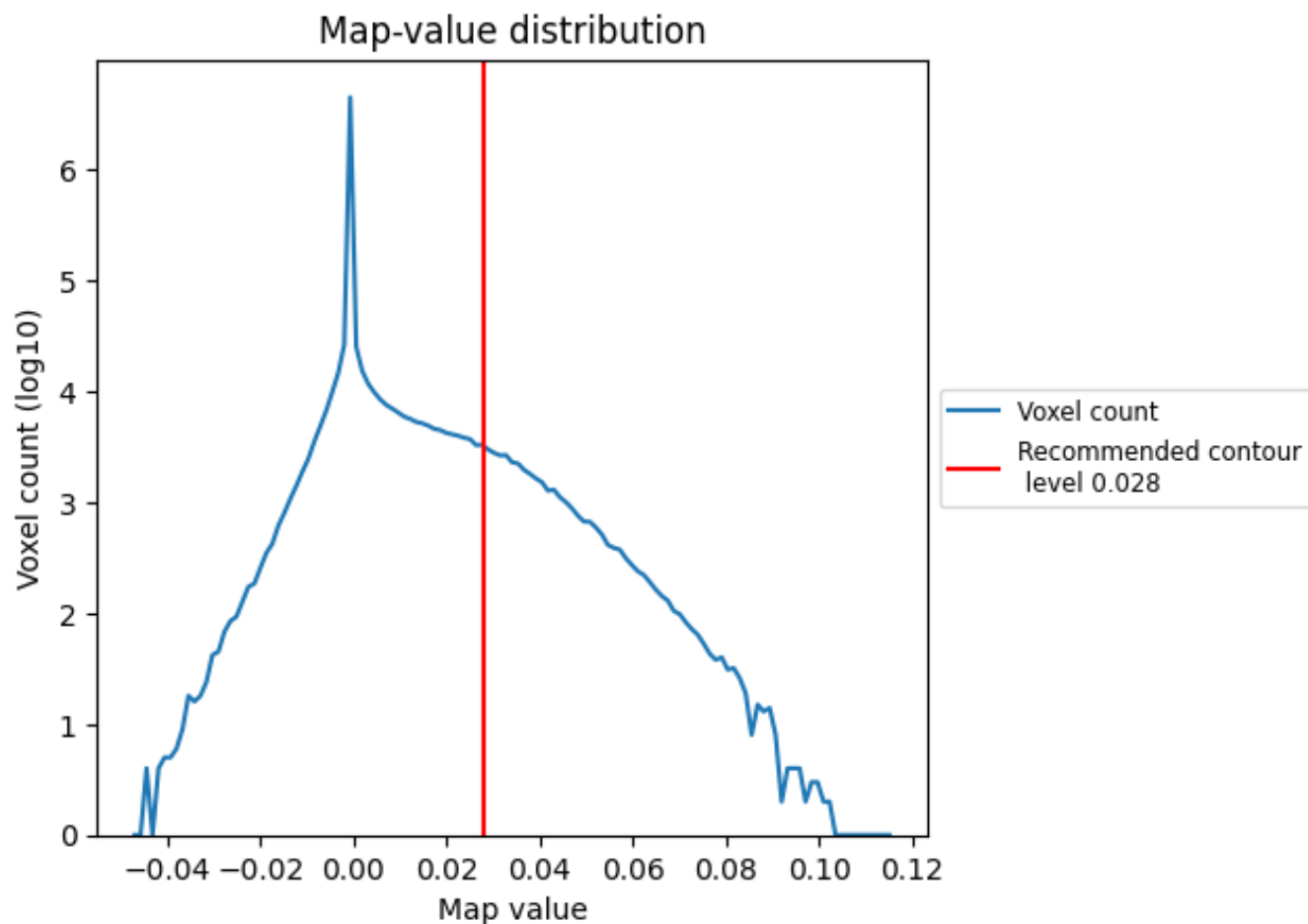
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

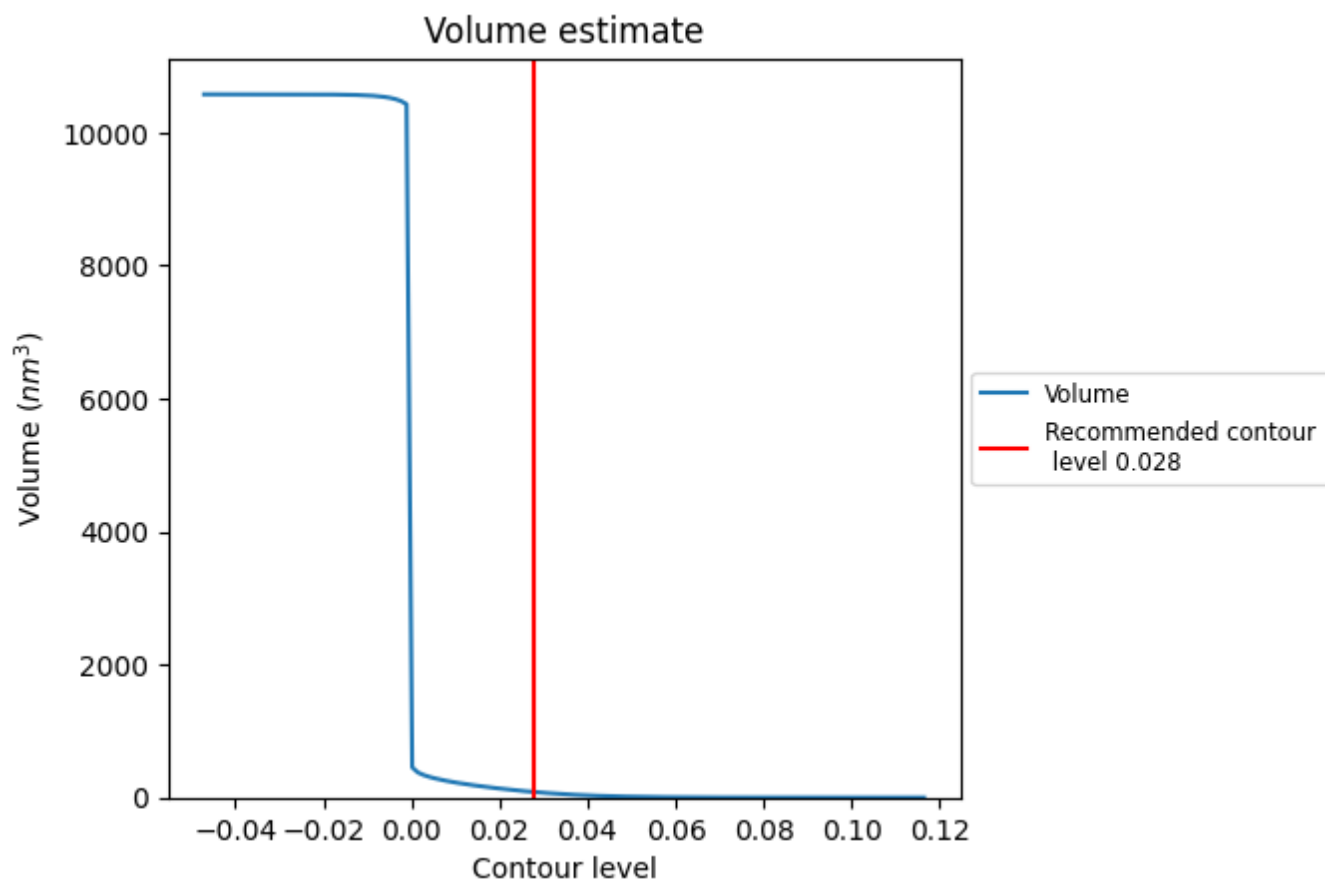
This section contains the results of statistical analysis of the map.

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.

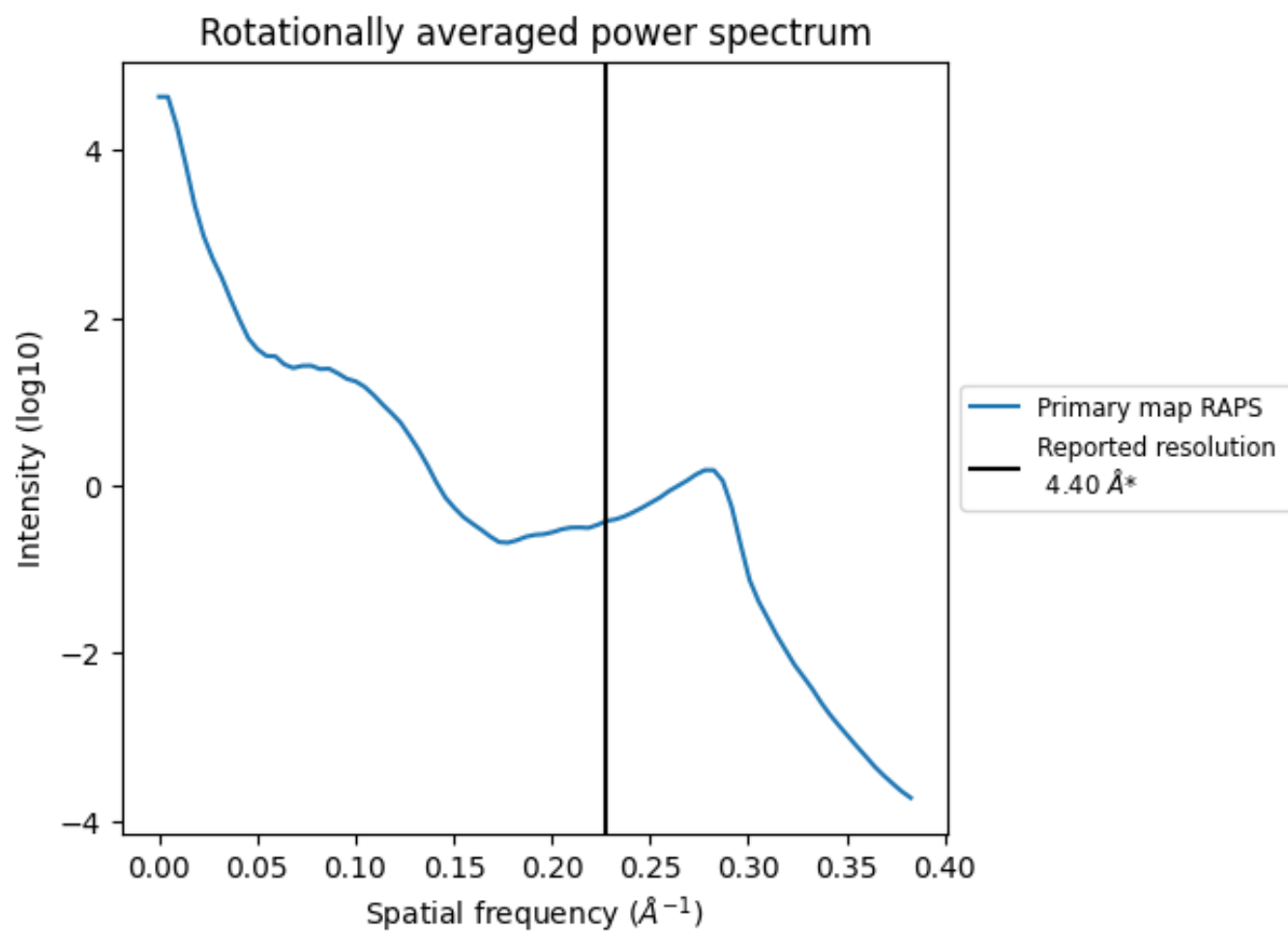
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 83 nm³; this corresponds to an approximate mass of 75 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 0.227 Å⁻¹

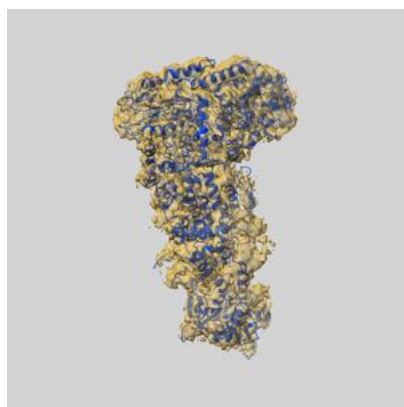
8 Fourier-Shell correlation

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

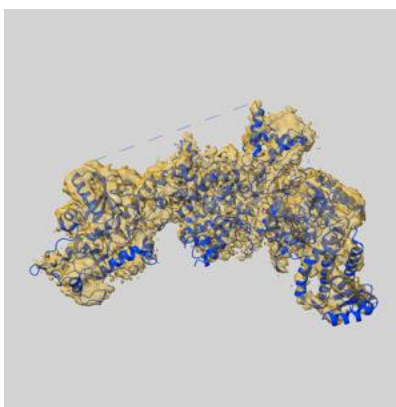
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-6904 and PDB model 5ZAK. Per-residue inclusion information can be found in section 3 on page 4.

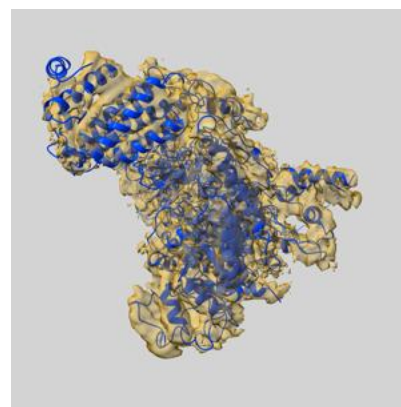
9.1 Map-model overlay [i](#)



X



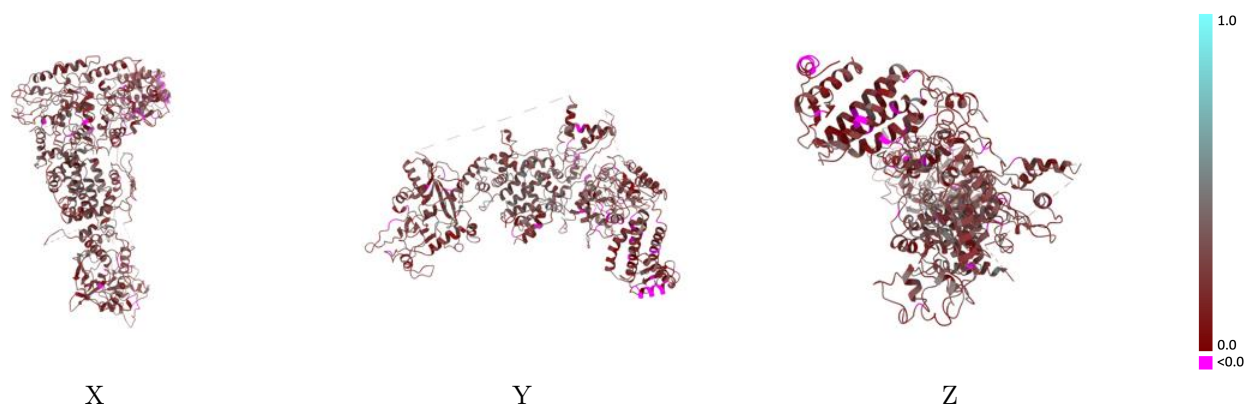
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.028 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](#)

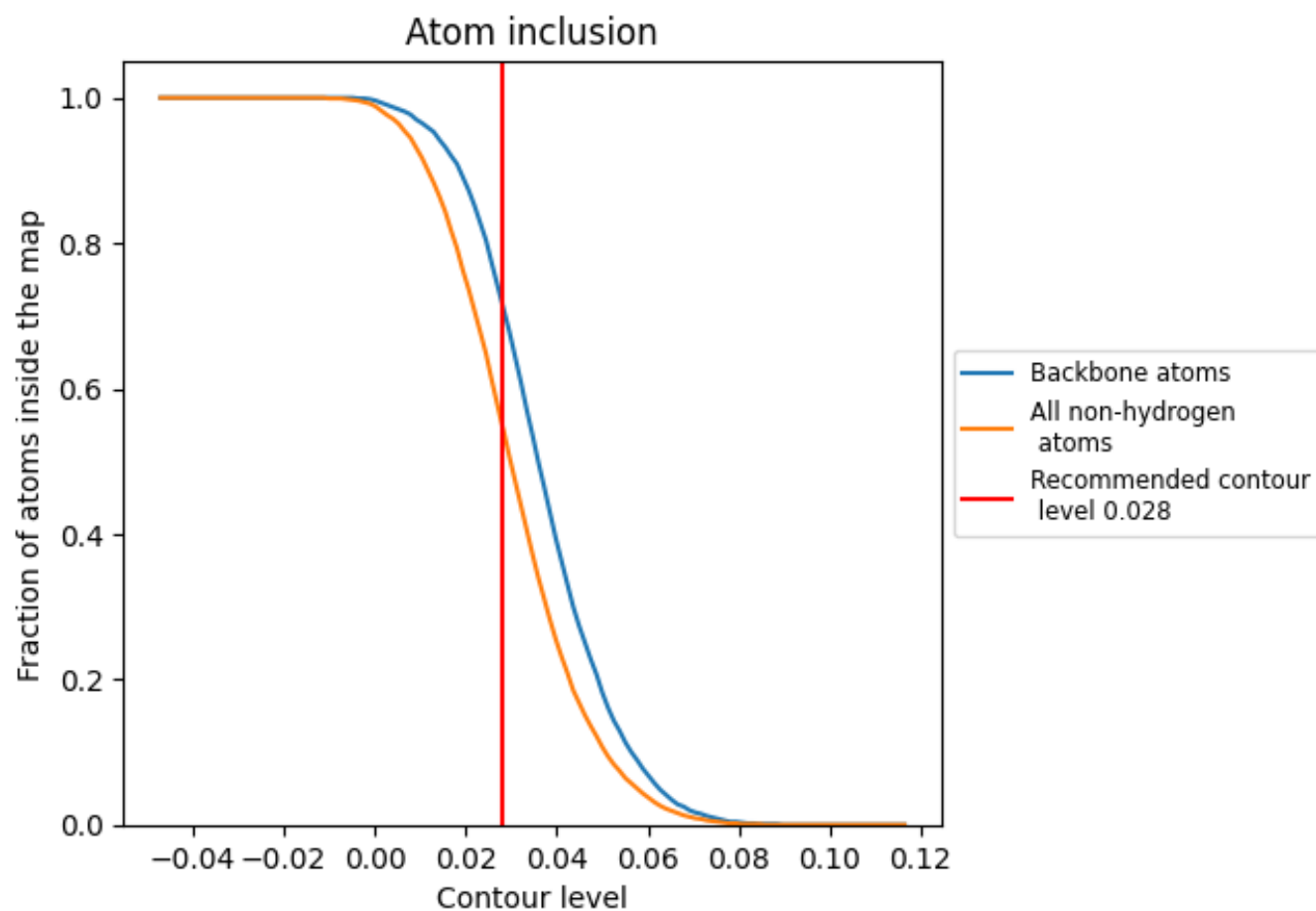


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 72% of all backbone atoms, 55% 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 (0.028) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5490	0.2500
A	0.5630	0.2540
B	0.2740	0.1680

