



Full wwPDB EM Validation Report ⓘ

Mar 23, 2026 – 08:08 PM UTC

PDB ID : 7DXK / pdb_00007dxk
EMDB ID : EMD-30912
Title : Human 128QHuntingtin-HAP40 complex structure
Authors : Guo, Q.; Fernandez-Busnadiego, R.
Deposited on : 2021-01-19
Resolution : 4.10 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

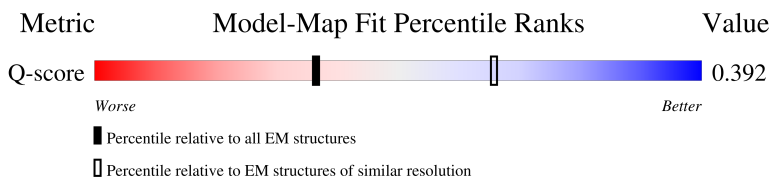
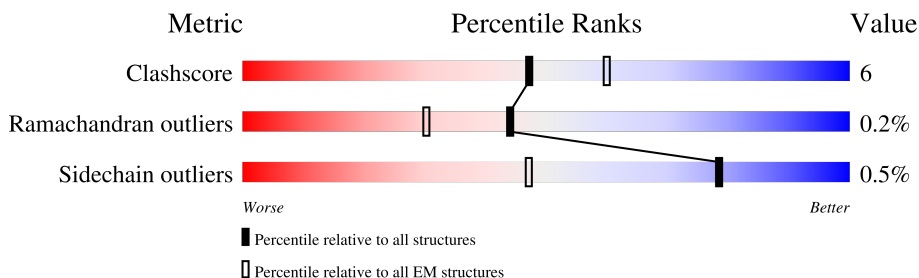
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.10 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	6458 (3.60 - 4.60)

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	3249	 5% 60% 12% 23%
2	B	371	 7% 60% 12% 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 20491 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 Huntingtin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	2353	Total	C	N	O	S	0	0
			18458	11839	3157	3343	119		

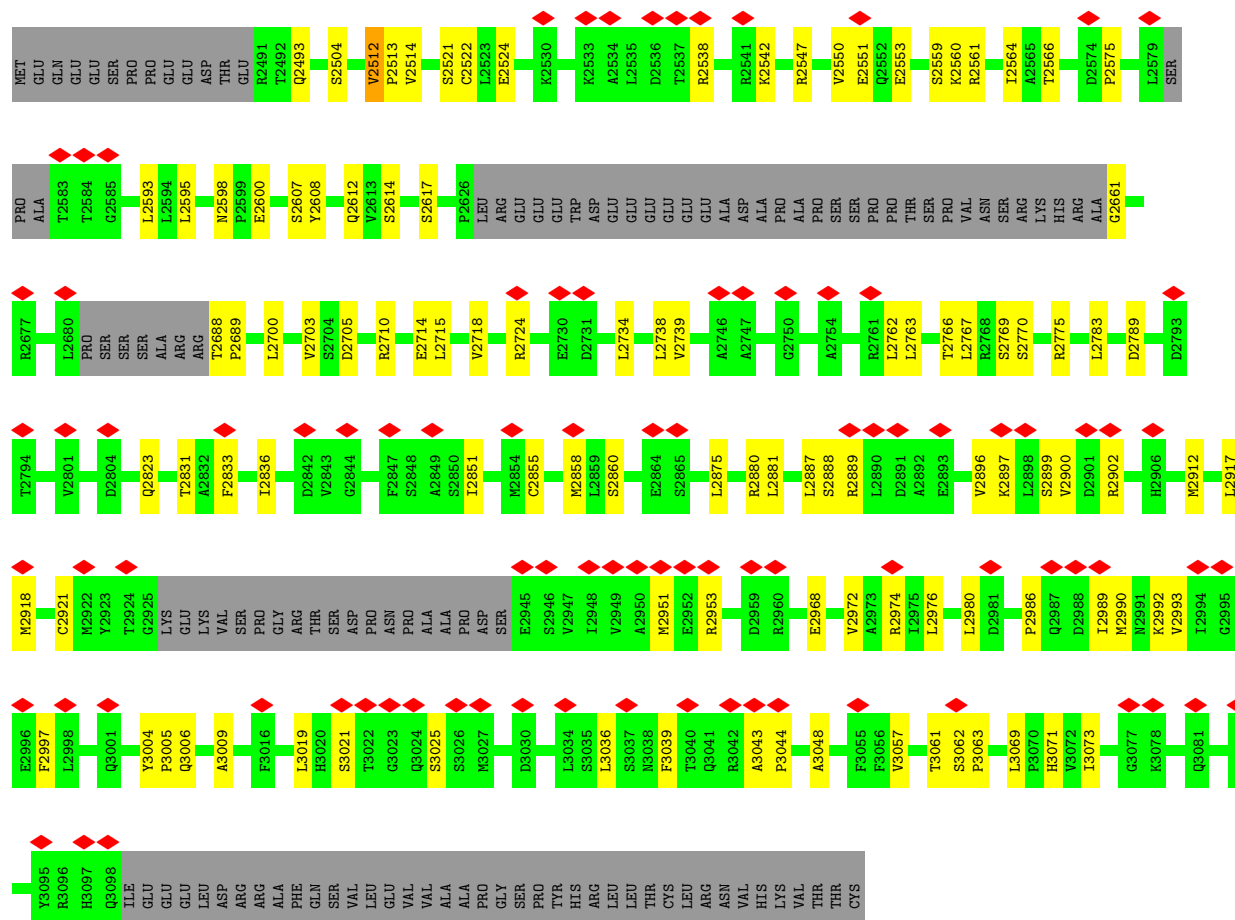
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1234	ARG	LYS	conflict	UNP P42858
A	1716	ASN	THR	variant	UNP P42858
A	2305	HIS	TYR	variant	UNP P42858

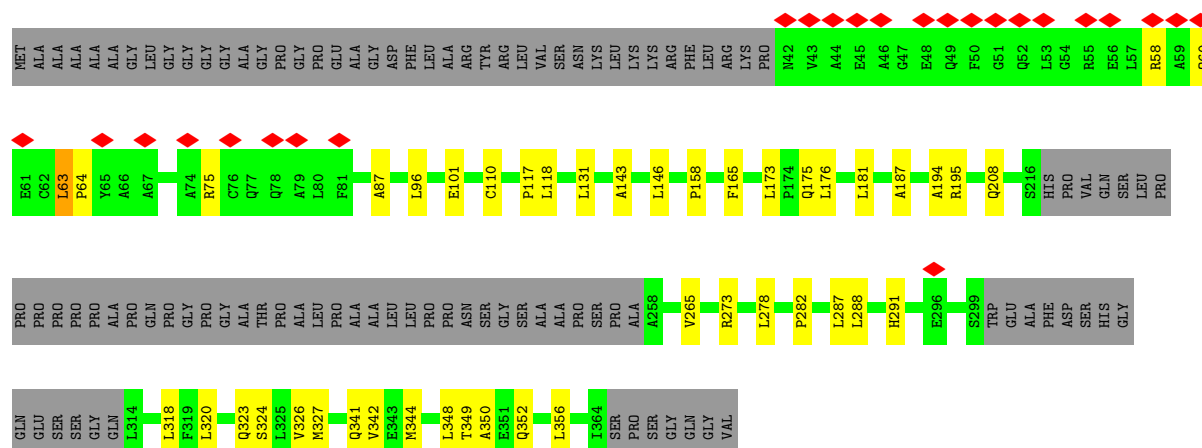
- Molecule 2 is a protein called 40-kDa huntingtin-associated protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	268	Total	C	N	O	S	0	0
			2033	1290	367	366	10		

L841	L849	L852	R853	T867	F878	V907	L911	L927	L928	L934	D939	D944	P945	Y946	Y947	X959	LEU	LEU	MET	HIS	GLU	THR	GLN	PRO	PRO	SER	HIS	PHE	SER	VAL	SER	THR	ILE	THR	R978	Y983	N996	T1014	R1015	L1025	C1026	L1027	P1033	



• Molecule 2: 40-kDa huntingtin-associated protein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	122992	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.412	Depositor
Minimum map value	-0.250	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	216.0, 216.0, 216.0	wwPDB
Map dimensions	160, 160, 160	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.35, 1.35, 1.35	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	0.30	0/18818	0.59	2/25566 (0.0%)
2	B	0.32	0/2067	0.59	1/2808 (0.0%)
All	All	0.30	0/20885	0.59	3/28374 (0.0%)

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	4
2	B	0	1
All	All	0	5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2514	VAL	N-CA-C	-8.54	104.81	113.10
2	B	63	LEU	N-CA-C	5.08	121.05	109.81
1	A	852	LEU	CA-CB-CG	5.01	133.83	116.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	276	HIS	Peptide
1	A	296	VAL	Peptide
1	A	3044	PRO	Peptide
1	A	3062	SER	Peptide
2	B	63	LEU	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	18458	0	18894	228	0
2	B	2033	0	2074	30	0
All	All	20491	0	20968	248	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 (248) 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:232:MET:HG3	1:A:273:ILE:HG12	1.70	0.74
1:A:2448:ARG:O	1:A:2452:LEU:HB2	1.88	0.74
1:A:2561:ARG:HE	2:B:195:ARG:HH12	1.36	0.71
1:A:1046:PRO:HG3	1:A:1060:VAL:HG23	1.76	0.65
1:A:2256:LEU:HD13	1:A:2266:PRO:HG2	1.80	0.63
1:A:2420:PRO:HB2	1:A:2427:ALA:H	1.64	0.61
1:A:275:GLN:HE21	1:A:320:LEU:HD21	1.66	0.61
1:A:2607:SER:O	1:A:2612:GLN:NE2	2.35	0.60
1:A:2547:ARG:NH1	1:A:2551:GLU:OE2	2.34	0.60
1:A:1147:ILE:HD11	1:A:1232:TYR:HB3	1.84	0.60
1:A:1461:LEU:HD13	1:A:1465:GLN:HA	1.83	0.59
1:A:3006:GLN:NE2	1:A:3073:ILE:O	2.34	0.59
2:B:58:ARG:HE	2:B:96:LEU:HD11	1.66	0.59
1:A:939:ASP:OD1	1:A:996:ASN:ND2	2.34	0.59
1:A:2521:SER:OG	1:A:2522:CYS:N	2.34	0.59
1:A:2271:LEU:HD21	1:A:2384:LEU:HD11	1.85	0.59
1:A:1963:PRO:HB2	2:B:320:LEU:HD23	1.85	0.59
1:A:1072:SER:OG	1:A:1073:ALA:N	2.36	0.58
1:A:150:VAL:O	1:A:154:LEU:HB3	2.04	0.58
1:A:252:ILE:HA	1:A:255:LEU:HD23	1.86	0.57
1:A:102:ILE:HG21	1:A:127:LEU:HD11	1.87	0.57
1:A:2769:SER:O	1:A:2775:ARG:NH2	2.38	0.56
1:A:1812:ASP:OD1	1:A:1812:ASP:N	2.38	0.56
1:A:256:LYS:HG2	1:A:306:LEU:HD12	1.87	0.56
1:A:2600:GLU:OE2	1:A:2823:GLN:NE2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2365:SER:O	1:A:2369:VAL:N	2.38	0.56
1:A:3057:VAL:HG13	1:A:3069:LEU:HD12	1.87	0.56
1:A:3021:SER:O	1:A:3025:SER:N	2.40	0.55
1:A:836:SER:OG	1:A:837:SER:N	2.39	0.55
1:A:1615:ALA:O	1:A:1617:GLN:NE2	2.40	0.55
1:A:2493:GLN:NE2	2:B:60:GLN:O	2.39	0.55
1:A:2860:SER:HA	1:A:2902:ARG:HD2	1.89	0.55
1:A:2208:ALA:O	1:A:2212:GLN:HB2	2.07	0.54
1:A:244:ILE:HD13	1:A:277:SER:HB2	1.90	0.54
2:B:282:PRO:HG2	2:B:287:LEU:HD12	1.89	0.54
1:A:1014:THR:OG1	1:A:1015:ARG:N	2.35	0.54
1:A:1823:THR:O	1:A:1886:ASN:ND2	2.39	0.54
1:A:138:ASP:OD1	1:A:138:ASP:N	2.39	0.54
1:A:804:ILE:HG21	1:A:841:LEU:HB3	1.89	0.54
1:A:1703:SER:HG	1:A:1705:THR:HG1	1.54	0.54
1:A:1659:MET:HG2	1:A:1710:LEU:HB3	1.89	0.54
1:A:2265:ILE:HD11	2:B:342:VAL:HA	1.89	0.53
1:A:260:PRO:HA	1:A:263:ARG:HE	1.74	0.53
1:A:1788:PHE:HB3	1:A:1791:ILE:HD12	1.89	0.53
1:A:1999:VAL:HG22	2:B:324:SER:HB2	1.90	0.53
1:A:2912:MET:HE1	1:A:2968:GLU:HA	1.90	0.53
1:A:206:LEU:HB2	1:A:247:LEU:HD21	1.90	0.53
1:A:2595:LEU:HB3	1:A:2880:ARG:HH11	1.73	0.53
1:A:181:ALA:HB1	1:A:185:ARG:HH12	1.74	0.53
1:A:1822:ILE:HA	1:A:1829:VAL:HG21	1.91	0.53
1:A:1435:THR:OG1	1:A:1436:THR:N	2.42	0.53
1:A:1960:LEU:HD13	1:A:1969:THR:HG21	1.90	0.53
1:A:1888:GLU:OE2	1:A:1891:ARG:NH1	2.40	0.53
2:B:101:GLU:HB2	2:B:117:PRO:HG2	1.90	0.53
2:B:181:LEU:HD13	2:B:265:VAL:HG11	1.91	0.52
1:A:1531:VAL:HA	1:A:1535:ILE:HD12	1.90	0.52
1:A:384:PRO:HG2	1:A:387:LEU:HB3	1.92	0.52
1:A:2980:LEU:HB3	1:A:2989:ILE:HG21	1.92	0.52
1:A:878:PHE:HB2	1:A:945:PRO:HB2	1.91	0.52
1:A:2259:HIS:HB3	1:A:2266:PRO:HG3	1.91	0.52
1:A:2444:GLU:OE2	1:A:2448:ARG:NH2	2.42	0.52
1:A:691:THR:O	1:A:744:GLN:NE2	2.43	0.52
1:A:1033:PRO:HG2	1:A:1036:ILE:HD12	1.91	0.52
1:A:2400:LEU:O	1:A:2426:THR:OG1	2.27	0.52
1:A:198:CYS:SG	1:A:199:ARG:N	2.83	0.52
1:A:2986:PRO:HG3	1:A:3019:LEU:HD11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2126:ASP:OD1	1:A:2126:ASP:N	2.40	0.51
1:A:826:VAL:HG11	1:A:849:VAL:HG21	1.92	0.51
1:A:2710:ARG:NH1	1:A:2714:GLU:OE2	2.44	0.51
1:A:151:ILE:HD13	1:A:162:LEU:HD13	1.91	0.51
2:B:158:PRO:HB2	2:B:194:ALA:HB2	1.93	0.51
1:A:2311:LEU:HD13	1:A:2388:LEU:HD13	1.92	0.51
1:A:682:ARG:HG2	1:A:730:LEU:HA	1.93	0.51
1:A:206:LEU:HD23	1:A:207:PRO:HD3	1.92	0.51
1:A:1419:ARG:HH21	1:A:1421:PHE:HB2	1.76	0.50
1:A:1597:ASP:OD1	1:A:1597:ASP:N	2.42	0.50
1:A:2614:SER:O	1:A:2617:SER:OG	2.26	0.50
1:A:2769:SER:OG	1:A:2770:SER:N	2.42	0.50
1:A:135:ALA:HA	1:A:177:ARG:HH22	1.77	0.50
1:A:2553:GLU:OE1	1:A:2974:ARG:NH1	2.43	0.50
1:A:2700:LEU:HA	1:A:2703:VAL:HG12	1.94	0.50
1:A:2151:GLU:HG2	1:A:2158:SER:HB2	1.94	0.49
1:A:853:ARG:HD3	1:A:907:VAL:HG22	1.93	0.49
1:A:2734:LEU:O	1:A:2738:LEU:HB2	2.12	0.49
1:A:2036:GLN:HE21	1:A:2055:LEU:HG	1.77	0.49
1:A:2598:ASN:OD1	1:A:2598:ASN:N	2.45	0.49
1:A:2457:ARG:NH2	1:A:2661:GLY:O	2.46	0.49
1:A:2899:SER:HB3	1:A:2917:LEU:HD23	1.93	0.49
1:A:2433:VAL:HG23	1:A:2436:LEU:HD12	1.95	0.49
1:A:2705:ASP:N	1:A:2705:ASP:OD2	2.45	0.49
1:A:983:TYR:OH	1:A:1272:GLN:O	2.25	0.48
1:A:2036:GLN:NE2	1:A:2056:ASP:OD2	2.46	0.48
2:B:208:GLN:OE1	2:B:273:ARG:NH2	2.45	0.48
1:A:1979:SER:O	1:A:2026:GLN:NE2	2.46	0.48
1:A:2315:ALA:HB2	1:A:2380:VAL:HG23	1.95	0.48
1:A:1128:ARG:HE	1:A:1131:VAL:HG21	1.78	0.48
1:A:1645:VAL:HB	1:A:1691:ARG:HG2	1.96	0.48
1:A:1747:LEU:HB3	1:A:1772:LEU:HD12	1.95	0.48
1:A:2547:ARG:HA	1:A:2550:VAL:HG12	1.96	0.48
1:A:365:ASP:HB3	1:A:368:VAL:HG12	1.94	0.48
1:A:778:ARG:O	1:A:780:ARG:NH2	2.47	0.48
1:A:2918:MET:HE2	1:A:2953:ARG:HG2	1.95	0.48
1:A:1796:THR:HG22	1:A:1839:VAL:HG23	1.96	0.48
2:B:143:ALA:HA	2:B:146:LEU:HD12	1.96	0.48
1:A:129:LEU:HD21	1:A:182:ALA:HB2	1.96	0.48
1:A:2119:VAL:HG13	1:A:2122:ILE:HD12	1.95	0.48
1:A:2993:VAL:O	1:A:2997:PHE:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1438:VAL:HG21	1:A:1485:GLU:HB3	1.96	0.48
1:A:2280:LEU:HD11	1:A:2566:THR:HG21	1.96	0.48
2:B:348:LEU:HD13	2:B:352:GLN:HB3	1.96	0.48
1:A:1132:PRO:O	1:A:1135:GLU:HB2	2.14	0.47
1:A:2397:ARG:NH1	1:A:2454:TRP:O	2.44	0.47
1:A:2881:LEU:HB3	1:A:2887:LEU:HD11	1.97	0.47
1:A:180:ARG:NH1	1:A:219:SER:OG	2.47	0.47
1:A:1649:LEU:HA	1:A:1652:MET:HB3	1.97	0.47
1:A:1684:THR:OG1	1:A:1685:GLU:N	2.41	0.47
1:A:2789:ASP:OD1	1:A:2789:ASP:N	2.43	0.47
1:A:2436:LEU:HD22	1:A:2442:PHE:HD1	1.79	0.47
1:A:2767:LEU:O	1:A:2775:ARG:NH1	2.47	0.47
1:A:827:ARG:HB2	1:A:867:THR:HG22	1.95	0.47
1:A:928:ILE:HD11	1:A:1027:LEU:HB3	1.97	0.47
1:A:944:ASP:HB2	1:A:947:VAL:HG23	1.95	0.47
1:A:1755:GLN:HB3	1:A:1758:VAL:HG22	1.97	0.47
1:A:2836:ILE:HD11	1:A:2851:ILE:HD13	1.97	0.47
1:A:2312:GLU:HG2	1:A:2316:VAL:HG21	1.97	0.47
1:A:2763:LEU:HA	1:A:2766:THR:HG22	1.95	0.47
1:A:1565:GLU:HA	1:A:1568:VAL:HG12	1.96	0.47
1:A:1308:VAL:HG11	1:A:1424:LEU:HD22	1.97	0.46
1:A:1900:ASP:HB2	1:A:1931:PRO:HB3	1.97	0.46
1:A:2951:MET:SD	1:A:2992:LYS:NZ	2.77	0.46
1:A:1844:TYR:OH	1:A:1911:HIS:ND1	2.43	0.46
1:A:667:GLY:O	1:A:679:HIS:ND1	2.48	0.46
1:A:1629:VAL:HA	1:A:1632:THR:HG22	1.98	0.46
1:A:1927:SER:O	1:A:1927:SER:OG	2.32	0.46
1:A:1570:MET:HE3	1:A:1570:MET:HB2	1.85	0.46
1:A:1646:ASP:OD2	1:A:1646:ASP:N	2.49	0.46
1:A:2972:VAL:O	1:A:2976:LEU:HB2	2.16	0.46
2:B:87:ALA:HB2	2:B:131:LEU:HD23	1.98	0.45
2:B:165:PHE:HB3	2:B:187:ALA:HB2	1.98	0.45
1:A:1364:GLN:NE2	1:A:1368:ASP:OD2	2.50	0.45
1:A:761:ARG:NH1	1:A:813:ASP:OD2	2.50	0.45
1:A:1742:GLN:HE22	1:A:1809:TYR:H	1.64	0.45
1:A:3004:TYR:HA	1:A:3005:PRO:HD3	1.86	0.45
1:A:93:LYS:HA	1:A:93:LYS:HD3	1.88	0.45
1:A:319:PRO:HG3	1:A:378:GLN:NE2	2.32	0.45
1:A:2176:VAL:HG21	1:A:2244:ASP:HB3	1.98	0.45
1:A:727:PHE:HZ	1:A:775:ILE:HD11	1.82	0.45
1:A:2538:ARG:O	1:A:2542:LYS:NZ	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2607:SER:OG	1:A:2608:TYR:N	2.50	0.45
2:B:349:THR:OG1	2:B:350:ALA:N	2.49	0.45
1:A:2356:CYS:SG	1:A:2357:GLU:N	2.90	0.45
1:A:1670:SER:OG	1:A:1771:GLU:OE2	2.35	0.44
1:A:2888:SER:OG	1:A:2889:ARG:N	2.50	0.44
2:B:101:GLU:HG2	2:B:118:LEU:HG	1.99	0.44
1:A:2171:ARG:HD3	1:A:2171:ARG:HA	1.86	0.44
1:A:2311:LEU:HD11	1:A:2370:LEU:HD21	1.99	0.44
1:A:3005:PRO:O	1:A:3009:ALA:N	2.49	0.44
1:A:752:TYR:HB2	1:A:764:THR:HG21	2.00	0.44
2:B:75:ARG:HE	2:B:75:ARG:HB3	1.68	0.44
2:B:288:LEU:HB2	2:B:291:HIS:HD2	1.82	0.44
1:A:3036:LEU:HA	1:A:3039:PHE:HB2	1.99	0.44
2:B:341:GLN:HG3	2:B:356:LEU:HD22	1.99	0.44
1:A:934:LEU:HD23	1:A:934:LEU:HA	1.80	0.44
1:A:1431:GLN:HG3	1:A:1440:LEU:HD21	1.98	0.44
1:A:2512:VAL:HG13	1:A:2513:PRO:HD3	2.00	0.44
1:A:2739:VAL:HG12	1:A:2762:LEU:HD21	2.00	0.44
1:A:1514:PRO:HA	1:A:1517:ILE:HD12	1.98	0.44
1:A:2875:LEU:HD11	1:A:2917:LEU:HB2	2.00	0.44
1:A:2221:LEU:HD23	1:A:2221:LEU:HA	1.85	0.44
1:A:772:ILE:HG12	1:A:800:LEU:HD12	1.99	0.43
1:A:907:VAL:O	1:A:911:LEU:HB2	2.18	0.43
1:A:1686:ASP:N	1:A:1686:ASP:OD1	2.50	0.43
1:A:169:GLU:OE1	1:A:178:SER:OG	2.33	0.43
1:A:1924:ILE:HD13	1:A:1969:THR:HG22	2.00	0.43
1:A:2993:VAL:O	1:A:2997:PHE:CB	2.66	0.43
1:A:1603:SER:HA	1:A:1606:ILE:HG22	2.00	0.43
1:A:1659:MET:HB2	1:A:1665:VAL:HG22	2.00	0.43
1:A:2524:GLU:HA	1:A:2575:PRO:HB2	2.00	0.43
1:A:2559:SER:OG	1:A:2560:LYS:N	2.52	0.43
2:B:344:MET:HE2	2:B:344:MET:HB3	1.78	0.43
1:A:3043:ALA:HB3	1:A:3048:ALA:HB2	2.01	0.43
1:A:698:VAL:O	1:A:708:LYS:NZ	2.50	0.43
1:A:1639:PRO:HB2	1:A:1643:ARG:HH21	1.83	0.43
1:A:927:LEU:HA	1:A:927:LEU:HD23	1.79	0.43
1:A:2833:PHE:HA	1:A:2836:ILE:HD12	1.99	0.43
1:A:3061:THR:HG23	1:A:3063:PRO:HD2	2.00	0.43
1:A:391:LEU:HD23	1:A:391:LEU:HA	1.91	0.43
1:A:2120:ASN:HA	1:A:2160:LEU:HD21	2.00	0.43
1:A:2381:PRO:HA	2:B:350:ALA:HB1	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1729:LYS:HA	1:A:1729:LYS:HD2	1.81	0.42
1:A:2783:LEU:HD13	1:A:2831:THR:HG22	2.01	0.42
1:A:1065:MET:O	1:A:1068:THR:OG1	2.32	0.42
1:A:2016:LEU:HD21	1:A:2058:PHE:HB2	2.02	0.42
1:A:159:LEU:HD21	1:A:192:LEU:HB3	2.01	0.42
1:A:2855:CYS:HA	1:A:2858:MET:HB2	2.00	0.42
1:A:695:ASN:N	1:A:695:ASN:OD1	2.50	0.42
1:A:1256:THR:O	1:A:1256:THR:OG1	2.34	0.42
1:A:1761:SER:OG	1:A:1762:GLU:N	2.51	0.42
1:A:2438:GLU:HB3	1:A:2441:VAL:HG12	2.01	0.42
1:A:96:VAL:HG13	1:A:142:VAL:HG11	2.01	0.42
1:A:748:ASP:O	1:A:751:ASN:HB2	2.19	0.42
1:A:1497:LEU:HA	1:A:1500:LEU:HD12	2.02	0.42
1:A:1489:ILE:HG12	1:A:1493:ILE:HD11	2.02	0.42
1:A:2378:SER:O	1:A:2378:SER:OG	2.31	0.42
1:A:2897:LYS:HA	1:A:2900:VAL:HG12	2.01	0.42
2:B:278:LEU:HD11	2:B:326:VAL:HG22	2.01	0.42
1:A:231:ILE:O	1:A:235:PHE:N	2.52	0.41
1:A:1025:LEU:HD23	1:A:1025:LEU:HA	1.83	0.41
1:A:1786:GLY:O	1:A:1789:ARG:NE	2.53	0.41
1:A:2419:LYS:HB3	1:A:2419:LYS:HE2	1.84	0.41
1:A:978:ARG:O	1:A:1313:LYS:NZ	2.40	0.41
1:A:1966:LEU:HD23	1:A:2000:LEU:HD23	2.02	0.41
1:A:1982:GLY:H	1:A:2026:GLN:NE2	2.19	0.41
2:B:318:LEU:HD12	2:B:318:LEU:HA	1.85	0.41
1:A:200:PRO:HA	1:A:203:VAL:HG12	2.03	0.41
1:A:1845:ARG:HE	1:A:1845:ARG:HB3	1.65	0.41
1:A:2002:ARG:HG2	2:B:327:MET:HE1	2.01	0.41
1:A:2504:SER:OG	2:B:110:CYS:SG	2.75	0.41
2:B:173:LEU:HD23	2:B:173:LEU:HA	1.92	0.41
1:A:1619:MET:HE2	1:A:1619:MET:HB2	1.80	0.41
1:A:1141:LEU:HD23	1:A:1141:LEU:HA	1.79	0.41
1:A:2688:THR:HA	1:A:2689:PRO:HD3	1.92	0.41
1:A:2896:VAL:HG22	1:A:2921:CYS:HB3	2.02	0.41
1:A:1524:MET:HE3	1:A:1524:MET:HB2	1.75	0.41
1:A:2118:LEU:HD11	1:A:3071:HIS:HB2	2.03	0.41
1:A:1299:ARG:HA	1:A:1299:ARG:HD2	1.90	0.41
1:A:2724:ARG:HE	1:A:2762:LEU:HD12	1.84	0.41
1:A:1027:LEU:HD23	1:A:1027:LEU:HA	1.89	0.41
1:A:1669:ILE:HG12	1:A:1743:LEU:HD11	2.02	0.41
1:A:1999:VAL:HG11	2:B:323:GLN:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2408:PRO:HG2	1:A:2411:VAL:HG23	2.03	0.41
1:A:1275:GLU:OE2	1:A:1310:GLN:NE2	2.49	0.41
1:A:1838:LEU:HD23	1:A:1838:LEU:HA	1.90	0.41
1:A:2986:PRO:O	1:A:2990:MET:HB2	2.21	0.41
1:A:1733:GLU:HG2	1:A:1790:ARG:HB3	2.02	0.40
1:A:120:LEU:HA	1:A:123:ILE:HG22	2.03	0.40
1:A:681:VAL:HG22	1:A:715:VAL:HG23	2.03	0.40
1:A:1737:SER:HB2	1:A:1798:LEU:HD21	2.03	0.40
1:A:1966:LEU:HD12	1:A:1966:LEU:HA	1.90	0.40
1:A:2388:LEU:HA	1:A:2391:ILE:HG22	2.03	0.40
1:A:2715:LEU:HA	1:A:2718:VAL:HG22	2.03	0.40
1:A:2448:ARG:HD3	1:A:2448:ARG:HA	1.83	0.40
1:A:2836:ILE:HD11	1:A:2851:ILE:HG21	2.03	0.40
1:A:1445:LEU:HA	1:A:1445:LEU:HD23	1.84	0.40
1:A:1058:CYS:N	1:A:1095:ALA:O	2.54	0.40
1:A:1976:ILE:HD12	1:A:1980:GLN:HB3	2.02	0.40
1:A:2268:SER:OG	2:B:341:GLN:NE2	2.32	0.40
2:B:175:GLN:HB3	2:B:176:LEU:HD12	2.03	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	2313/3249 (71%)	2187 (95%)	123 (5%)	3 (0%)	48	81
2	B	262/371 (71%)	249 (95%)	12 (5%)	1 (0%)	30	65
All	All	2575/3620 (71%)	2436 (95%)	135 (5%)	4 (0%)	44	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	SER
1	A	1046	PRO
2	B	64	PRO
1	A	2266	PRO

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	2075/2874 (72%)	2064 (100%)	11 (0%)	81	82
2	B	202/274 (74%)	202 (100%)	0	100	100
All	All	2277/3148 (72%)	2266 (100%)	11 (0%)	78	82

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	746	VAL
1	A	748	ASP
1	A	1645	VAL
1	A	1702	ILE
1	A	1768	TYR
1	A	2122	ILE
1	A	2303	LEU
1	A	2385	THR
1	A	2512	VAL
1	A	2564	ILE
1	A	2593	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	ASN
1	A	148	ASN
1	A	158	ASN
1	A	204	ASN

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Mol	Chain	Res	Type
1	A	221	GLN
1	A	237	ASN
1	A	240	ASN
1	A	275	GLN
1	A	281	GLN
1	A	289	ASN
1	A	367	ASN
1	A	378	GLN
1	A	906	ASN
1	A	1140	HIS
1	A	1352	HIS
1	A	1589	GLN
1	A	1693	GLN
1	A	1708	ASN
1	A	1742	GLN
1	A	1765	HIS
1	A	1825	HIS
1	A	1928	HIS
1	A	1959	ASN
1	A	2026	GLN
1	A	2036	GLN
1	A	2040	GLN
1	A	2156	GLN
1	A	2223	GLN
1	A	2367	GLN
1	A	2377	ASN
1	A	2495	ASN
1	A	2529	ASN
1	A	2555	GLN
1	A	2571	GLN
1	A	2622	ASN
1	A	2665	HIS
1	A	2727	HIS
1	A	2809	ASN
1	A	3097	HIS
2	B	137	GLN
2	B	157	GLN
2	B	352	GLN

5.3.3 RNA ⓘ

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.

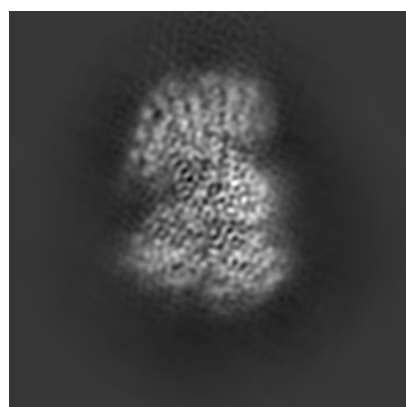
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30912. These allow visual inspection of the internal detail of the map and identification of artifacts.

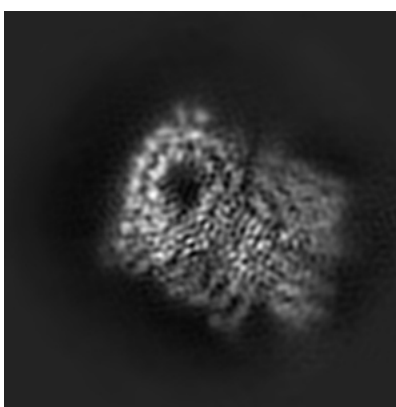
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

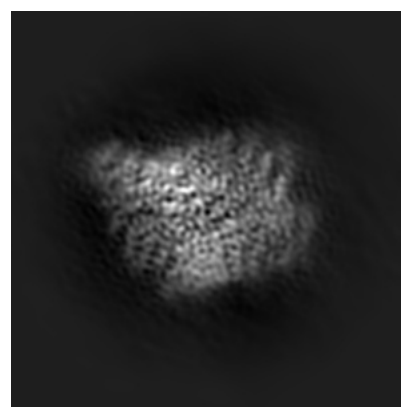
6.1.1 Primary map



X



Y

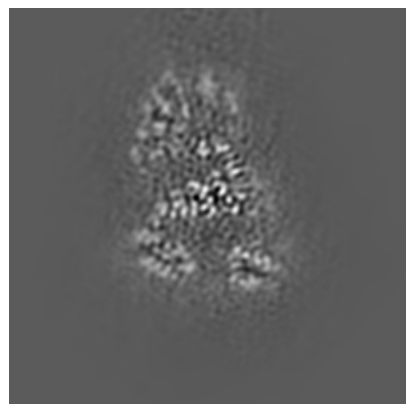


Z

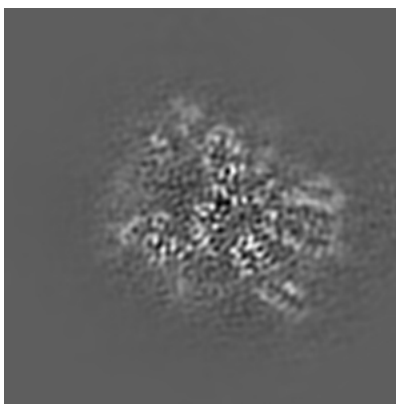
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

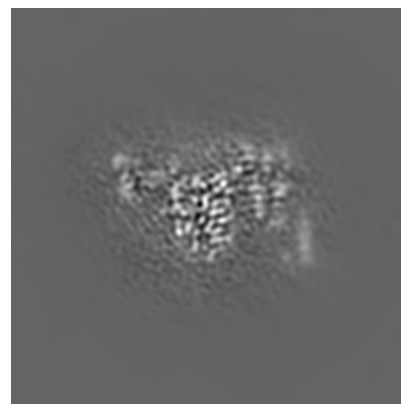
6.2.1 Primary map



X Index: 80



Y Index: 80

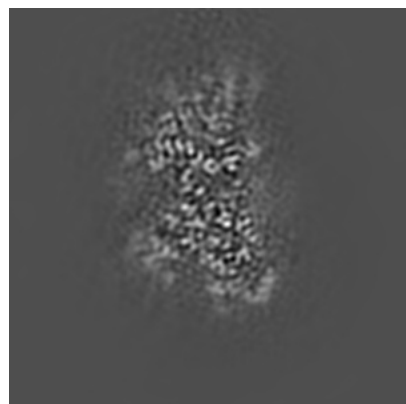


Z Index: 80

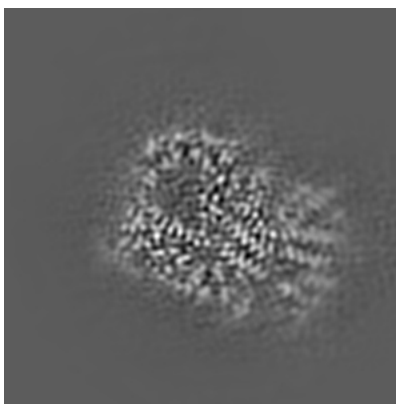
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

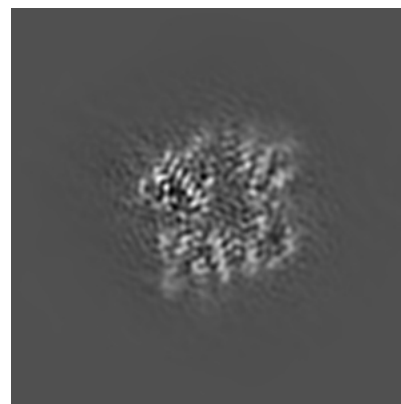
6.3.1 Primary map



X Index: 67



Y Index: 88

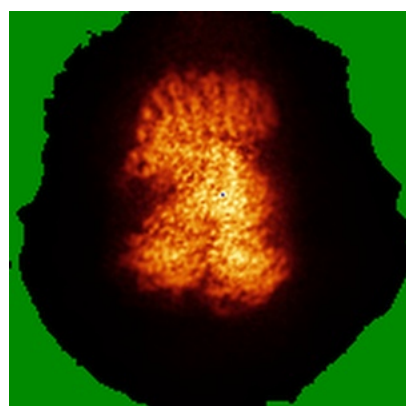


Z Index: 62

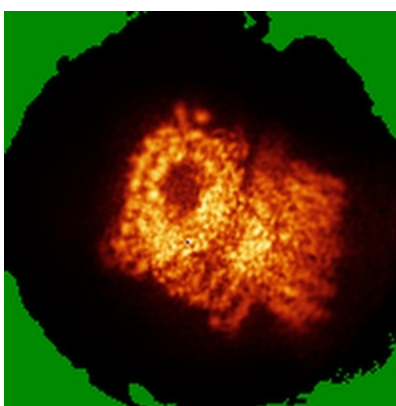
The images above show the largest variance slices of the map in three orthogonal directions.

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

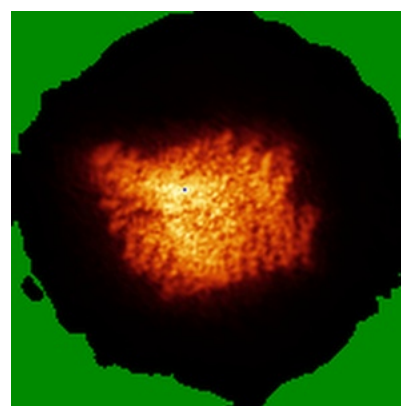
6.4.1 Primary map



X



Y

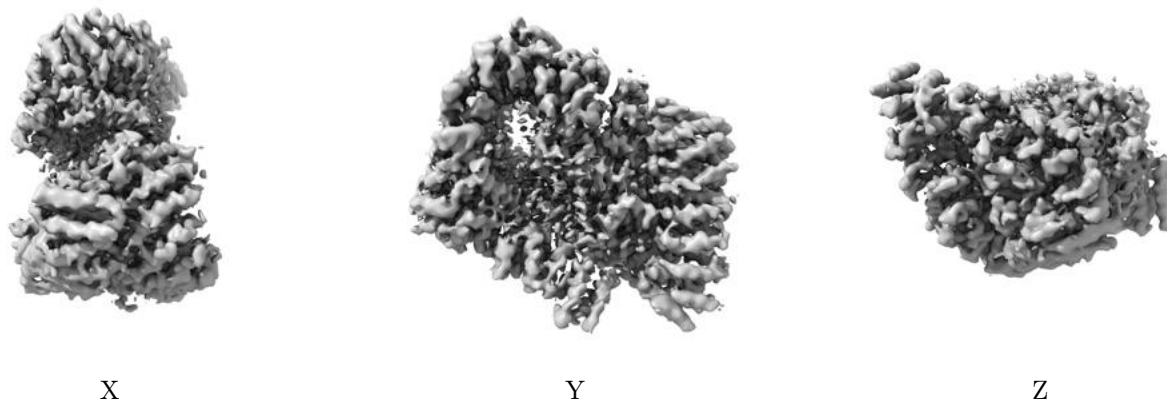


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

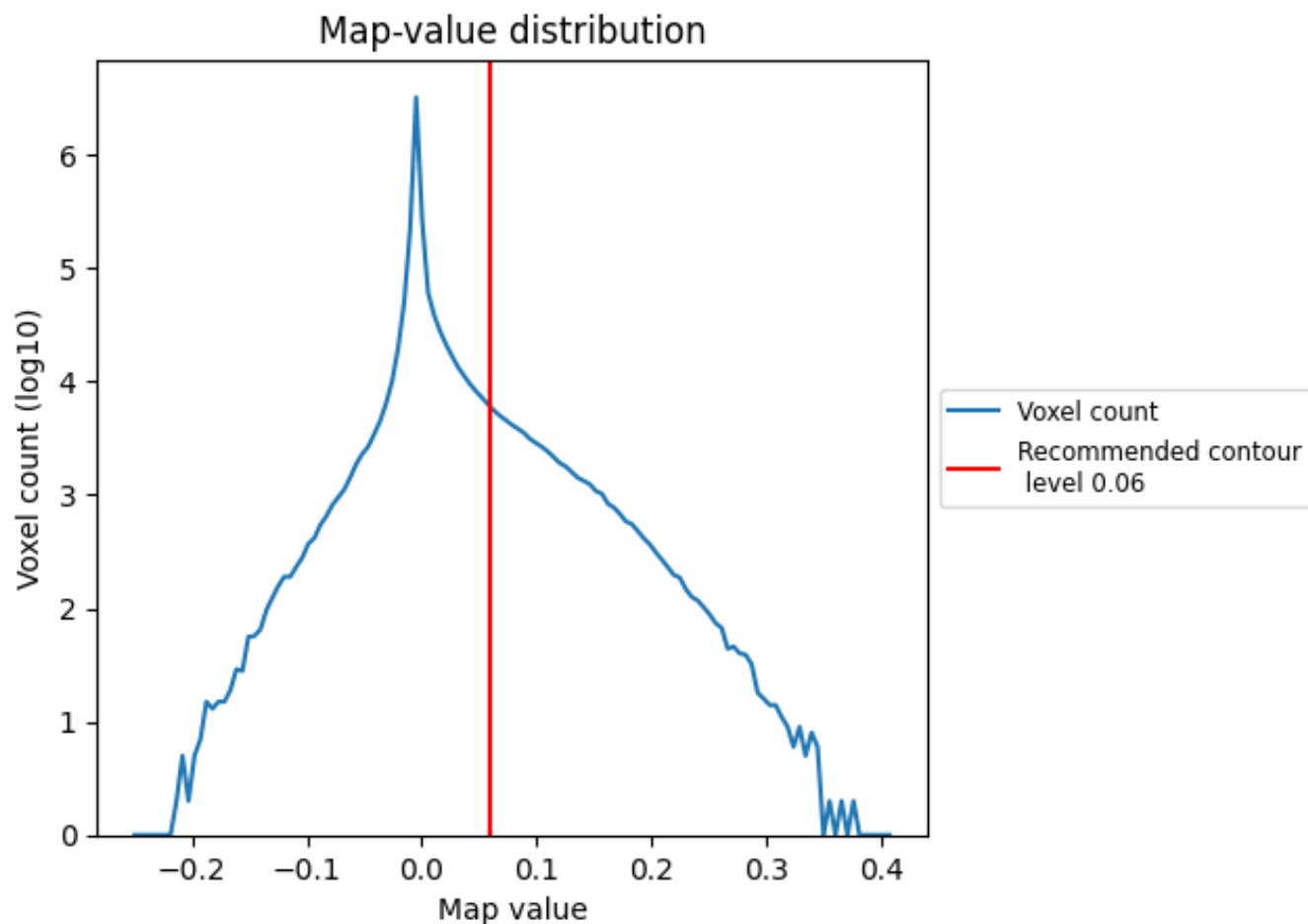
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

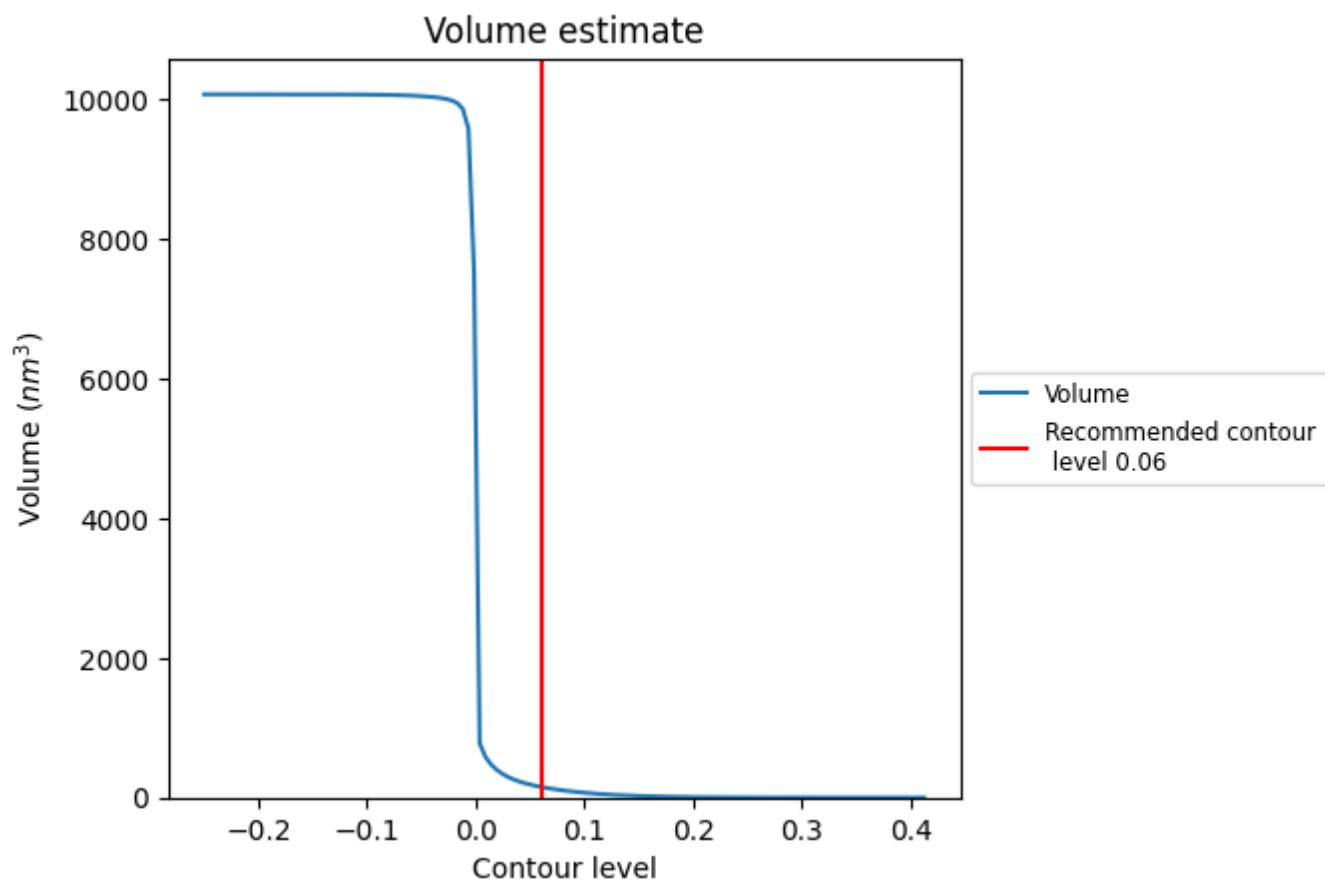
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

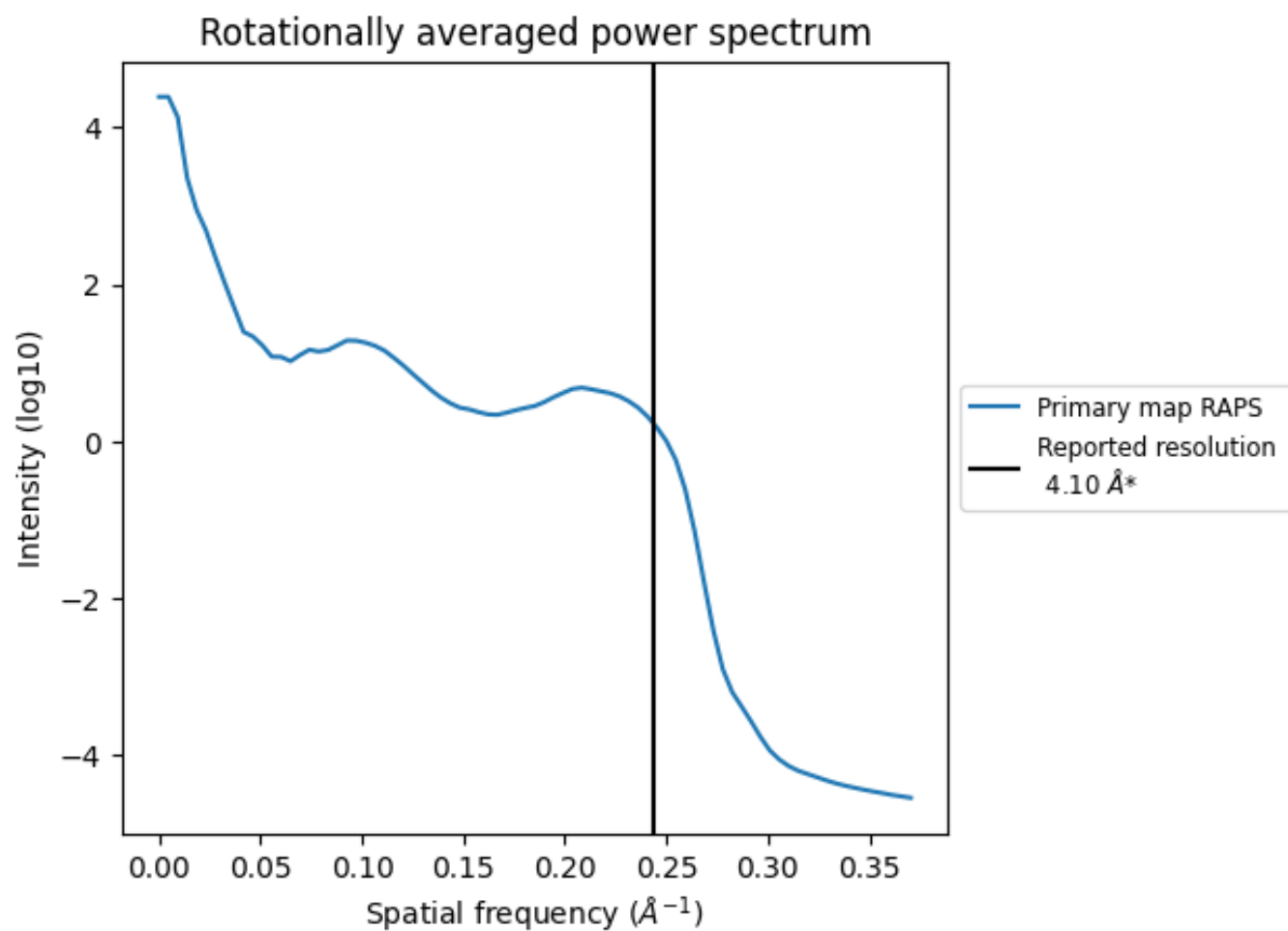
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 151 nm³; this corresponds to an approximate mass of 137 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 0.244 Å⁻¹

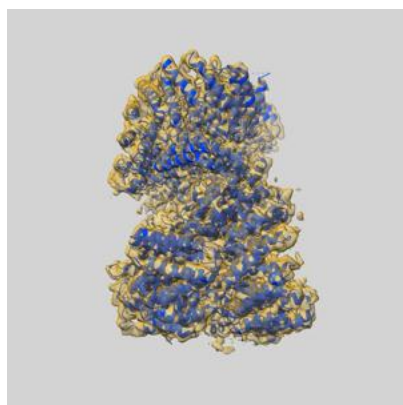
8 Fourier-Shell correlation ⓘ

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

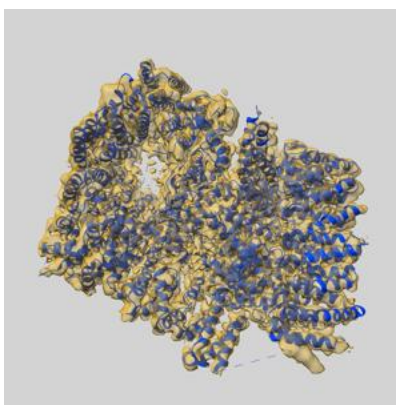
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30912 and PDB model 7DXK. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

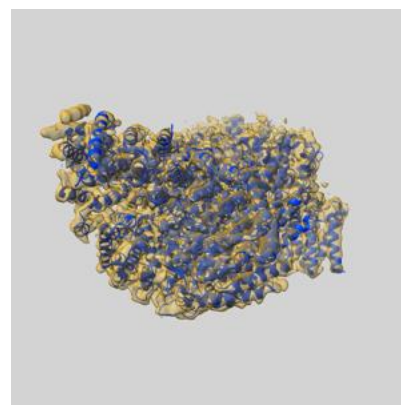
9.1 Map-model overlay [i](#)



X



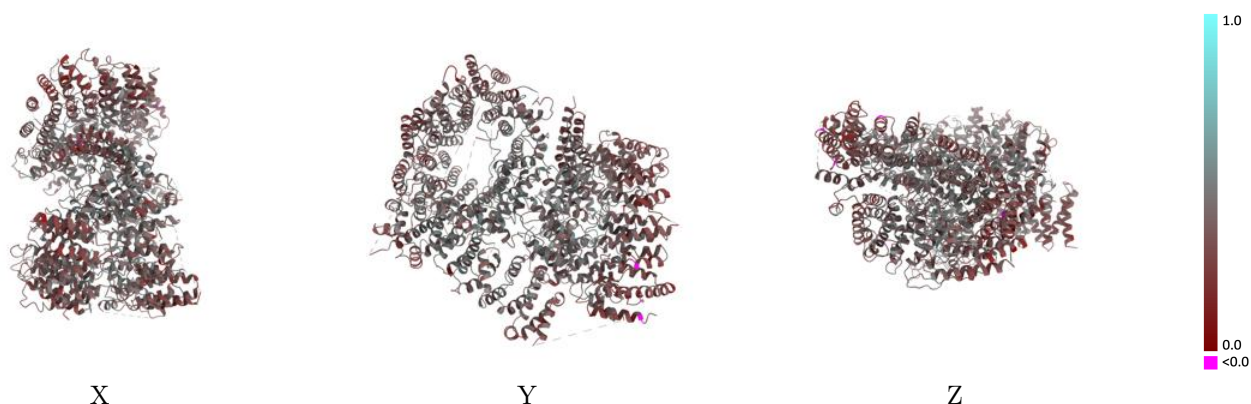
Y



Z

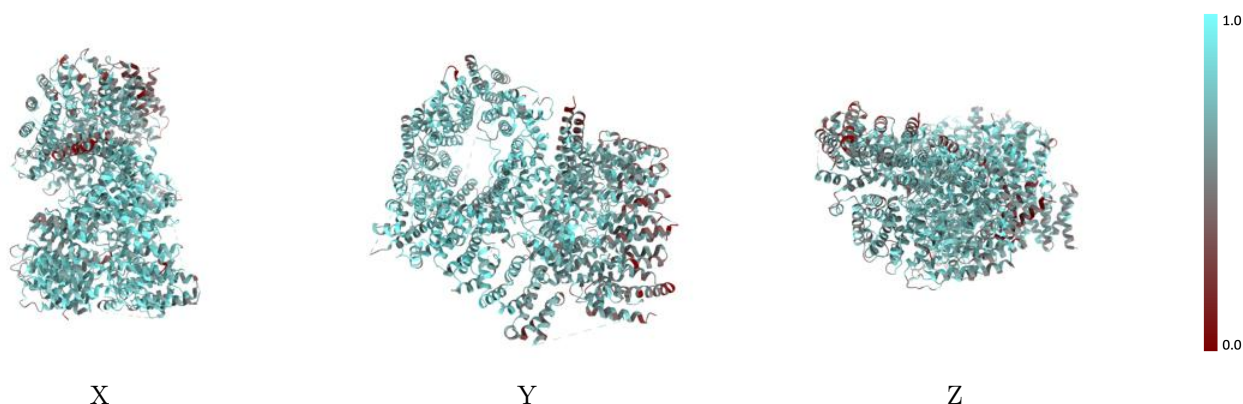
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



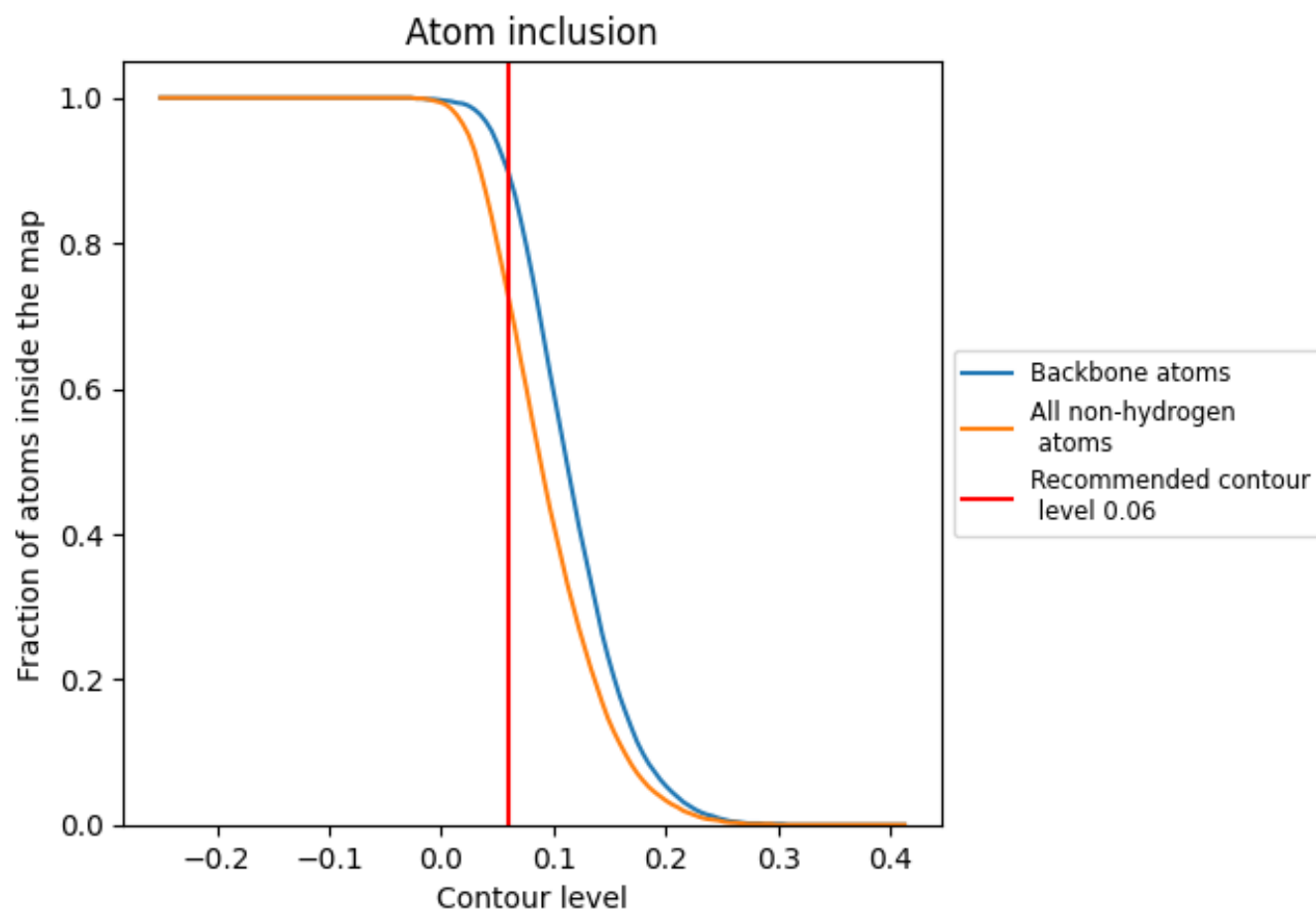
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.06).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 72% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7250	0.3920
A	0.7250	0.3890
B	0.7180	0.4190

