



wwPDB EM Validation Summary Report ⓘ

Aug 28, 2025 – 01:21 pm BST

PDB ID : 9HMS / pdb_00009hms
EMDB ID : EMD-52297
Title : Cryo-EM structure of human separase bound to SCC1 (310-550 aa) and SA2
Authors : Yu, J.; Schmidt, S.; Botto, M.; Boland, A.
Deposited on : 2024-12-09
Resolution : 3.40 Å (reported)
Based on initial models : 4PJU, 7NJ1

This is a wwPDB EM Validation Summary 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.dev126
MolProbity : 4-5-2 with Phenix2.0rc1
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

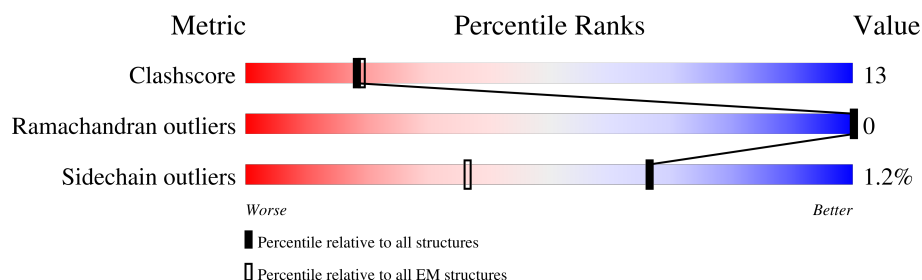
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.40 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1271	
2	B	241	
3	C	2172	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16909 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 Cohesin subunit SA-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	933	7178	4547	1230	1360	41	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1232	SER	-	expression tag	UNP Q8N3U4
A	1233	SER	-	expression tag	UNP Q8N3U4
A	1234	LEU	-	expression tag	UNP Q8N3U4
A	1235	ALA	-	expression tag	UNP Q8N3U4
A	1236	GLU	-	expression tag	UNP Q8N3U4
A	1237	GLU	-	expression tag	UNP Q8N3U4
A	1238	ASN	-	expression tag	UNP Q8N3U4
A	1239	LEU	-	expression tag	UNP Q8N3U4
A	1240	TYR	-	expression tag	UNP Q8N3U4
A	1241	PHE	-	expression tag	UNP Q8N3U4
A	1242	GLN	-	expression tag	UNP Q8N3U4
A	1243	SER	-	expression tag	UNP Q8N3U4
A	1244	TRP	-	expression tag	UNP Q8N3U4
A	1245	SER	-	expression tag	UNP Q8N3U4
A	1246	HIS	-	expression tag	UNP Q8N3U4
A	1247	PRO	-	expression tag	UNP Q8N3U4
A	1248	GLN	-	expression tag	UNP Q8N3U4
A	1249	PHE	-	expression tag	UNP Q8N3U4
A	1250	GLU	-	expression tag	UNP Q8N3U4
A	1251	LYS	-	expression tag	UNP Q8N3U4
A	1252	GLY	-	expression tag	UNP Q8N3U4
A	1253	GLY	-	expression tag	UNP Q8N3U4
A	1254	GLY	-	expression tag	UNP Q8N3U4
A	1255	SER	-	expression tag	UNP Q8N3U4
A	1256	GLY	-	expression tag	UNP Q8N3U4
A	1257	GLY	-	expression tag	UNP Q8N3U4
A	1258	GLY	-	expression tag	UNP Q8N3U4
A	1259	SER	-	expression tag	UNP Q8N3U4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1260	GLY	-	expression tag	UNP Q8N3U4
A	1261	GLY	-	expression tag	UNP Q8N3U4
A	1262	GLY	-	expression tag	UNP Q8N3U4
A	1263	SER	-	expression tag	UNP Q8N3U4
A	1264	TRP	-	expression tag	UNP Q8N3U4
A	1265	SER	-	expression tag	UNP Q8N3U4
A	1266	HIS	-	expression tag	UNP Q8N3U4
A	1267	PRO	-	expression tag	UNP Q8N3U4
A	1268	GLN	-	expression tag	UNP Q8N3U4
A	1269	PHE	-	expression tag	UNP Q8N3U4
A	1270	GLU	-	expression tag	UNP Q8N3U4
A	1271	LYS	-	expression tag	UNP Q8N3U4

- Molecule 2 is a protein called Double-strand-break repair protein rad21 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	107	Total	C	N	O	S	0	0
			869	552	145	168	4		

- Molecule 3 is a protein called Separin.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	1141	Total	C	N	O	S	0	0
			8861	5652	1536	1632	41		

There are 113 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-54	MET	-	initiating methionine	UNP Q14674
C	-53	ASP	-	expression tag	UNP Q14674
C	-52	TYR	-	expression tag	UNP Q14674
C	-51	LYS	-	expression tag	UNP Q14674
C	-50	ASP	-	expression tag	UNP Q14674
C	-49	HIS	-	expression tag	UNP Q14674
C	-48	ASP	-	expression tag	UNP Q14674
C	-47	GLY	-	expression tag	UNP Q14674
C	-46	ASP	-	expression tag	UNP Q14674
C	-45	TYR	-	expression tag	UNP Q14674
C	-44	LYS	-	expression tag	UNP Q14674
C	-43	ASP	-	expression tag	UNP Q14674
C	-42	HIS	-	expression tag	UNP Q14674
C	-41	ASP	-	expression tag	UNP Q14674

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-40	ILE	-	expression tag	UNP Q14674
C	-39	ASP	-	expression tag	UNP Q14674
C	-38	TYR	-	expression tag	UNP Q14674
C	-37	LYS	-	expression tag	UNP Q14674
C	-36	ASP	-	expression tag	UNP Q14674
C	-35	ASP	-	expression tag	UNP Q14674
C	-34	ASP	-	expression tag	UNP Q14674
C	-33	ASP	-	expression tag	UNP Q14674
C	-32	LYS	-	expression tag	UNP Q14674
C	-31	SER	-	expression tag	UNP Q14674
C	-30	GLY	-	expression tag	UNP Q14674
C	-29	PRO	-	expression tag	UNP Q14674
C	-28	GLY	-	expression tag	UNP Q14674
C	-27	GLY	-	expression tag	UNP Q14674
C	-26	SER	-	expression tag	UNP Q14674
C	-25	GLY	-	expression tag	UNP Q14674
C	-24	GLY	-	expression tag	UNP Q14674
C	-23	SER	-	expression tag	UNP Q14674
C	-22	GLY	-	expression tag	UNP Q14674
C	-21	GLY	-	expression tag	UNP Q14674
C	-20	GLY	-	expression tag	UNP Q14674
C	-19	SER	-	expression tag	UNP Q14674
C	-18	GLY	-	expression tag	UNP Q14674
C	-17	GLY	-	expression tag	UNP Q14674
C	-16	GLY	-	expression tag	UNP Q14674
C	-15	SER	-	expression tag	UNP Q14674
C	-14	GLY	-	expression tag	UNP Q14674
C	-13	GLU	-	expression tag	UNP Q14674
C	-12	ASN	-	expression tag	UNP Q14674
C	-11	LEU	-	expression tag	UNP Q14674
C	-10	TYR	-	expression tag	UNP Q14674
C	-9	PHE	-	expression tag	UNP Q14674
C	-8	GLN	-	expression tag	UNP Q14674
C	-7	GLY	-	expression tag	UNP Q14674
C	-6	GLY	-	expression tag	UNP Q14674
C	-5	GLY	-	expression tag	UNP Q14674
C	-4	SER	-	expression tag	UNP Q14674
C	-3	GLY	-	expression tag	UNP Q14674
C	-2	GLY	-	expression tag	UNP Q14674
C	-1	SER	-	expression tag	UNP Q14674
C	0	GLY	-	expression tag	UNP Q14674
C	25	ASP	ALA	conflict	UNP Q14674

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Chain	Residue	Modelled	Actual	Comment	Reference
C	116	VAL	ALA	conflict	UNP Q14674
C	1372	SER	ARG	conflict	UNP Q14674
C	1525	GLY	-	linker	UNP Q14674
C	1526	LEU	-	linker	UNP Q14674
C	1527	GLU	-	linker	UNP Q14674
C	1528	VAL	-	linker	UNP Q14674
C	1529	LEU	-	linker	UNP Q14674
C	1530	PHE	-	linker	UNP Q14674
C	1531	GLN	-	linker	UNP Q14674
C	1532	GLY	-	linker	UNP Q14674
C	1533	PRO	-	linker	UNP Q14674
C	1534	GLY	-	linker	UNP Q14674
C	1535	SER	-	linker	UNP Q14674
C	1536	GLY	-	linker	UNP Q14674
C	1561	GLN	ARG	conflict	UNP Q14674
C	2029	SER	CYS	conflict	UNP Q14674
C	2037	HIS	ARG	conflict	UNP Q14674
C	2121	SER	-	expression tag	UNP Q14674
C	2122	SER	-	expression tag	UNP Q14674
C	2123	LEU	-	expression tag	UNP Q14674
C	2124	ALA	-	expression tag	UNP Q14674
C	2125	GLU	-	expression tag	UNP Q14674
C	2126	GLU	-	expression tag	UNP Q14674
C	2127	ASN	-	expression tag	UNP Q14674
C	2128	LEU	-	expression tag	UNP Q14674
C	2129	TYR	-	expression tag	UNP Q14674
C	2130	PHE	-	expression tag	UNP Q14674
C	2131	GLN	-	expression tag	UNP Q14674
C	2132	SER	-	expression tag	UNP Q14674
C	2133	TRP	-	expression tag	UNP Q14674
C	2134	SER	-	expression tag	UNP Q14674
C	2135	HIS	-	expression tag	UNP Q14674
C	2136	PRO	-	expression tag	UNP Q14674
C	2137	GLN	-	expression tag	UNP Q14674
C	2138	PHE	-	expression tag	UNP Q14674
C	2139	GLU	-	expression tag	UNP Q14674
C	2140	LYS	-	expression tag	UNP Q14674
C	2141	GLY	-	expression tag	UNP Q14674
C	2142	GLY	-	expression tag	UNP Q14674
C	2143	GLY	-	expression tag	UNP Q14674
C	2144	SER	-	expression tag	UNP Q14674
C	2145	GLY	-	expression tag	UNP Q14674

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2146	GLY	-	expression tag	UNP Q14674
C	2147	GLY	-	expression tag	UNP Q14674
C	2148	SER	-	expression tag	UNP Q14674
C	2149	GLY	-	expression tag	UNP Q14674
C	2150	GLY	-	expression tag	UNP Q14674
C	2151	GLY	-	expression tag	UNP Q14674
C	2152	SER	-	expression tag	UNP Q14674
C	2153	TRP	-	expression tag	UNP Q14674
C	2154	SER	-	expression tag	UNP Q14674
C	2155	HIS	-	expression tag	UNP Q14674
C	2156	PRO	-	expression tag	UNP Q14674
C	2157	GLN	-	expression tag	UNP Q14674
C	2158	PHE	-	expression tag	UNP Q14674
C	2159	GLU	-	expression tag	UNP Q14674
C	2160	LYS	-	expression tag	UNP Q14674

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	C	1	Total	Zn	0
			1	1	





PHE	G2028	M1907	R1821	A1714
GLU	S2029	P1908	L1826	G1722
LYS	S2030	S1909	Q1827	N1723
GLY	L2034	L1910	W1831	T1724
GLY	G2042	L1913	K1832	L1727
SER	G2042	P1914	Y1833	T1728
GLY	L2056	V1915	T1837	R1729
GLY	F2057	T1916	K1840	L1730
GLY	L2058	R1917	I1841	P1735
GLY	L2061	F1921	M1842	P1736
GLY	W2062	F1923	L1843	V1737
SER	D2068	Y1927	G1847	S1738
SER	I2069	S1928	P1851	V1739
TRP	Y2072	I1929	I1854	Q1740
HIS	PRO	I1930	L1857	I1741
GLN	P2088	K1931	A1858	M1746
PHE	Y2092	E1932	Y1859	K1747
GLU	V2093	Y1933	L1860	L1750
LYS		G1934	G1861	F1757
		A1935	C1862	I1760
		L1939	P1863	M1767
	R2097	L1952	T1864	W1776
	P2100	E1962	Q1865	R1780
	R2101	E1972	P1866	L1781
	L2102	W1975	E1867	A1782
	K2103	V1978	Q1870	L1783
	L2104	V1979	E1871	D1784
	L2105	P1983	L1872	M1787
	I2106	V1988	N1874	I1791
	G2107	Y1998	E1875	
	P2110	G2006	R1879	V1798
	Y2113	A2007	L1880	W1802
	P2116	R2008	Q1881	K1803
	R2120	F2009	G1882	G1804
SER	SER	L2010	L1883	
SER	LEU	R2011	S1887	L1807
LEU	ALA	G2012	N1888	P1808
ALA	GLU	V2015	S1889	S1809
GLU	GLU	L2016	H1890	S1810
GLU	ASN	W2023	L1891	E1811
ASN	LEU	A2024	V1892	P1812
LEU	TYR	L2025	L1893	G1814
TYR	PHE	F2027	K1901	P1815
GLN	GLN		L1902	E1818
SER	SER		P1903	
TRP	TRP		S1906	
SER	SER			
HIS	HIS			
PRO	PRO			
GLN	GLN			
PHE	PHE			
GLU	GLU			
LYS	LYS			

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	177878	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	130000	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	22.198	Depositor
Minimum map value	-0.111	Depositor
Average map value	-0.022	Depositor
Map value standard deviation	0.291	Depositor
Recommended contour level	4.4	Depositor
Map size (Å)	462.0288, 462.0288, 462.0288	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.9024, 0.9024, 0.9024	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	0.52	0/7292	0.75	2/9865 (0.0%)
2	B	0.53	0/882	0.74	1/1188 (0.1%)
3	C	0.46	0/9036	0.67	3/12290 (0.0%)
All	All	0.49	0/17210	0.71	6/23343 (0.0%)

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	730	TYR	N-CA-C	-6.66	105.31	113.50
2	B	395	PRO	N-CA-CB	6.60	110.26	103.00
3	C	2027	PHE	N-CA-CB	-5.93	101.77	110.49
1	A	479	TYR	CB-CA-C	5.35	119.95	110.85
3	C	1798	VAL	N-CA-C	5.35	116.01	110.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7178	0	6844	174	0
2	B	869	0	878	26	0
3	C	8861	0	9021	245	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
All	All	16909	0	16743	434	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 13.

The worst 5 of 434 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:358:LYS:HE3	1:A:400:SER:HB3	1.64	0.78
3:C:1682:LEU:HD13	3:C:1690:PRO:HA	1.67	0.75
3:C:894:SER:HB3	3:C:897:TRP:HD1	1.50	0.75
1:A:970:ILE:HD11	1:A:1006:PHE:HB3	1.69	0.73
3:C:1880:LEU:HD12	3:C:1883:LEU:HD11	1.71	0.71

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	925/1271 (73%)	872 (94%)	53 (6%)	0	100	100
2	B	101/241 (42%)	94 (93%)	7 (7%)	0	100	100
3	C	1126/2172 (52%)	1053 (94%)	73 (6%)	0	100	100
All	All	2152/3684 (58%)	2019 (94%)	133 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	725/1136 (64%)	717 (99%)	8 (1%)	70	81
2	B	100/224 (45%)	97 (97%)	3 (3%)	36	61
3	C	969/1814 (53%)	958 (99%)	11 (1%)	70	81
All	All	1794/3174 (56%)	1772 (99%)	22 (1%)	66	80

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	941	ASP
3	C	1737	VAL
3	C	1597	TYR
3	C	1739	VAL
1	A	813	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 39 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	1208	HIS
3	C	1785	HIS
3	C	1237	GLN
3	C	1637	HIS
3	C	1953	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 1 ligands modelled in this entry, 1 is 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

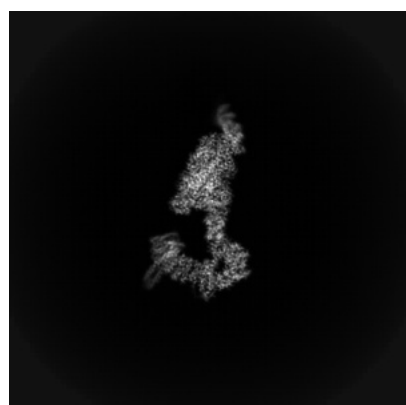
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-52297. These allow visual inspection of the internal detail of the map and identification of artifacts.

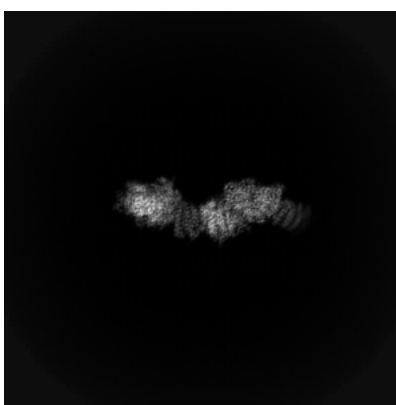
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

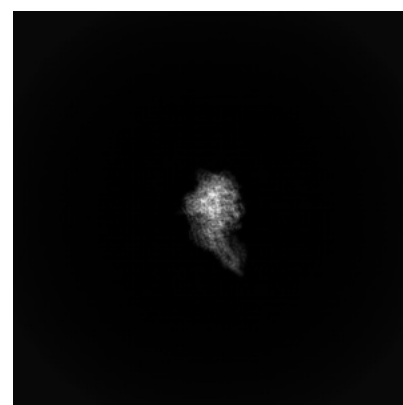
6.1.1 Primary map



X



Y

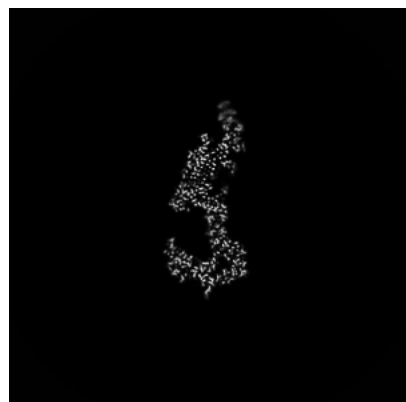


Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 256



Y Index: 256

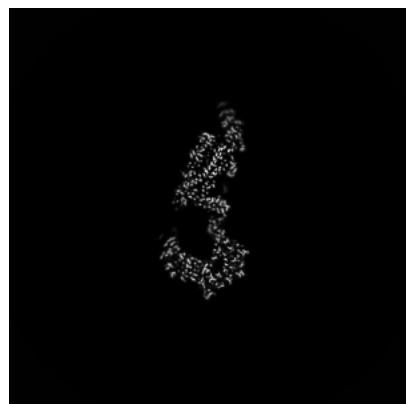


Z Index: 256

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: 261



Y Index: 259



Z Index: 175

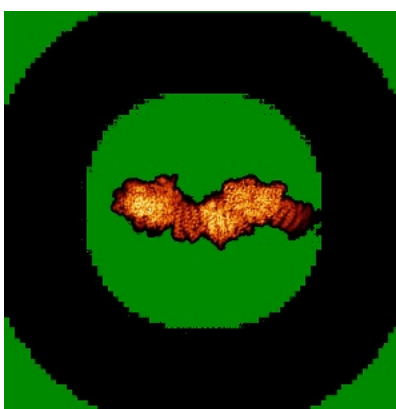
The images above show the largest variance slices of the map in three orthogonal directions.

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

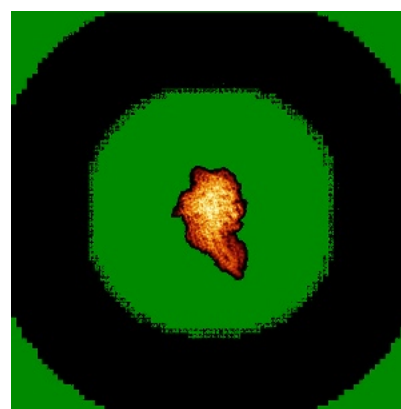
6.4.1 Primary map



X



Y

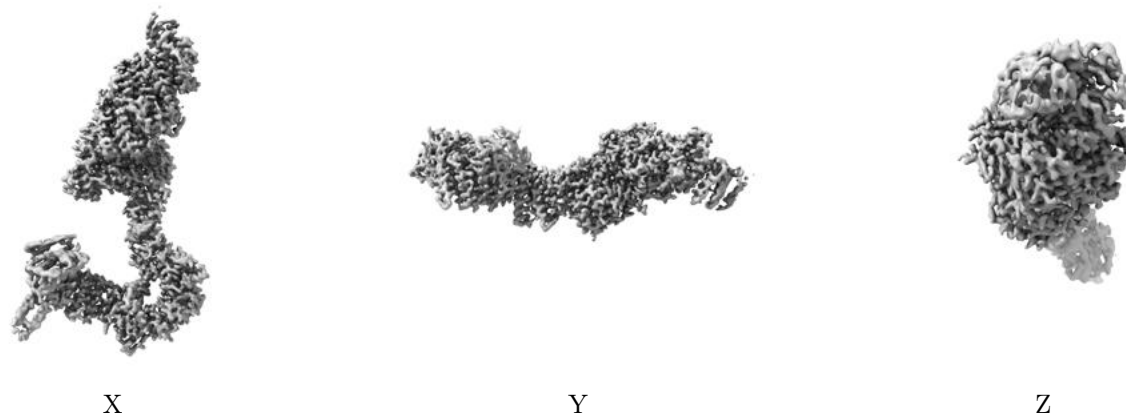


Z

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

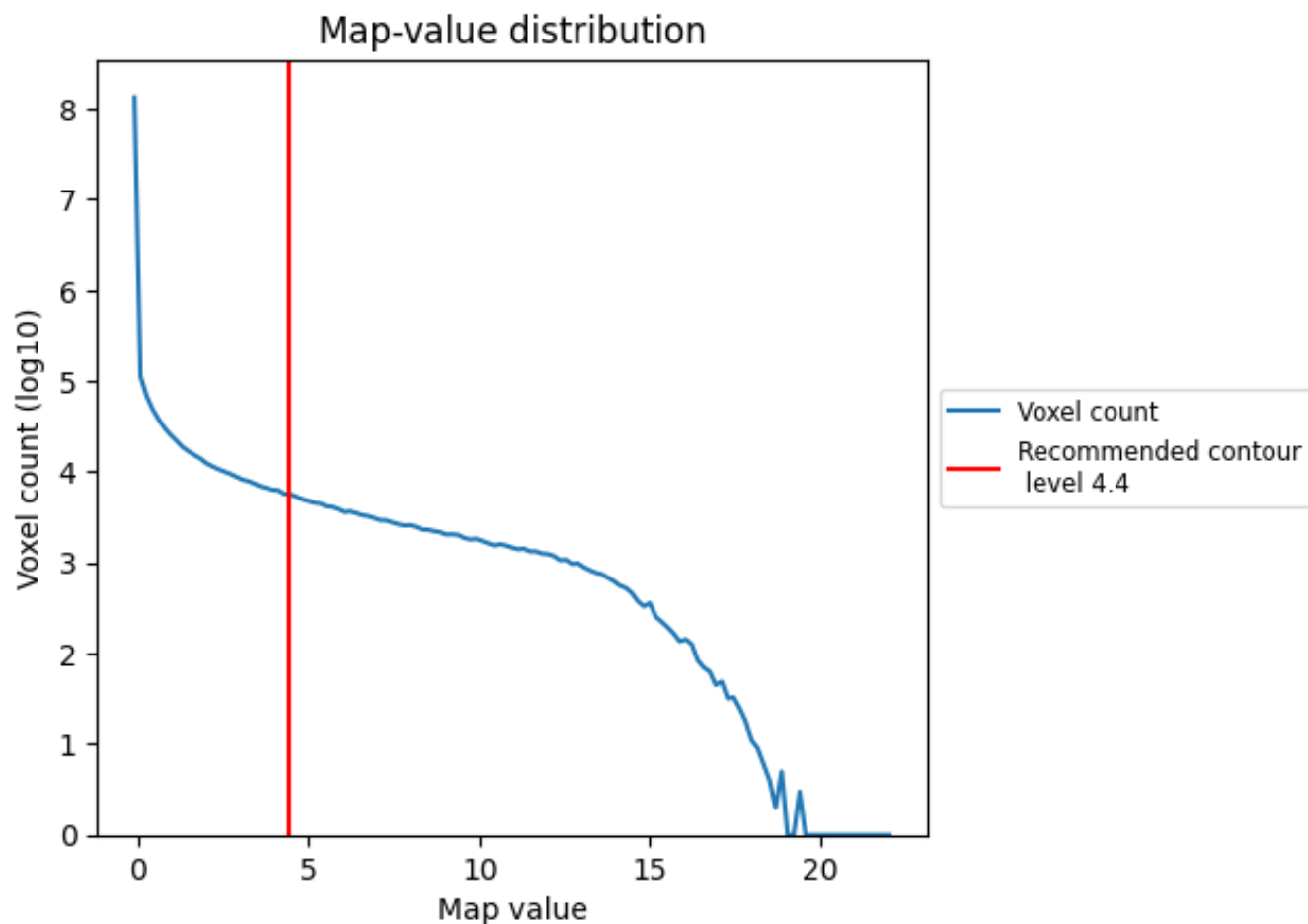
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

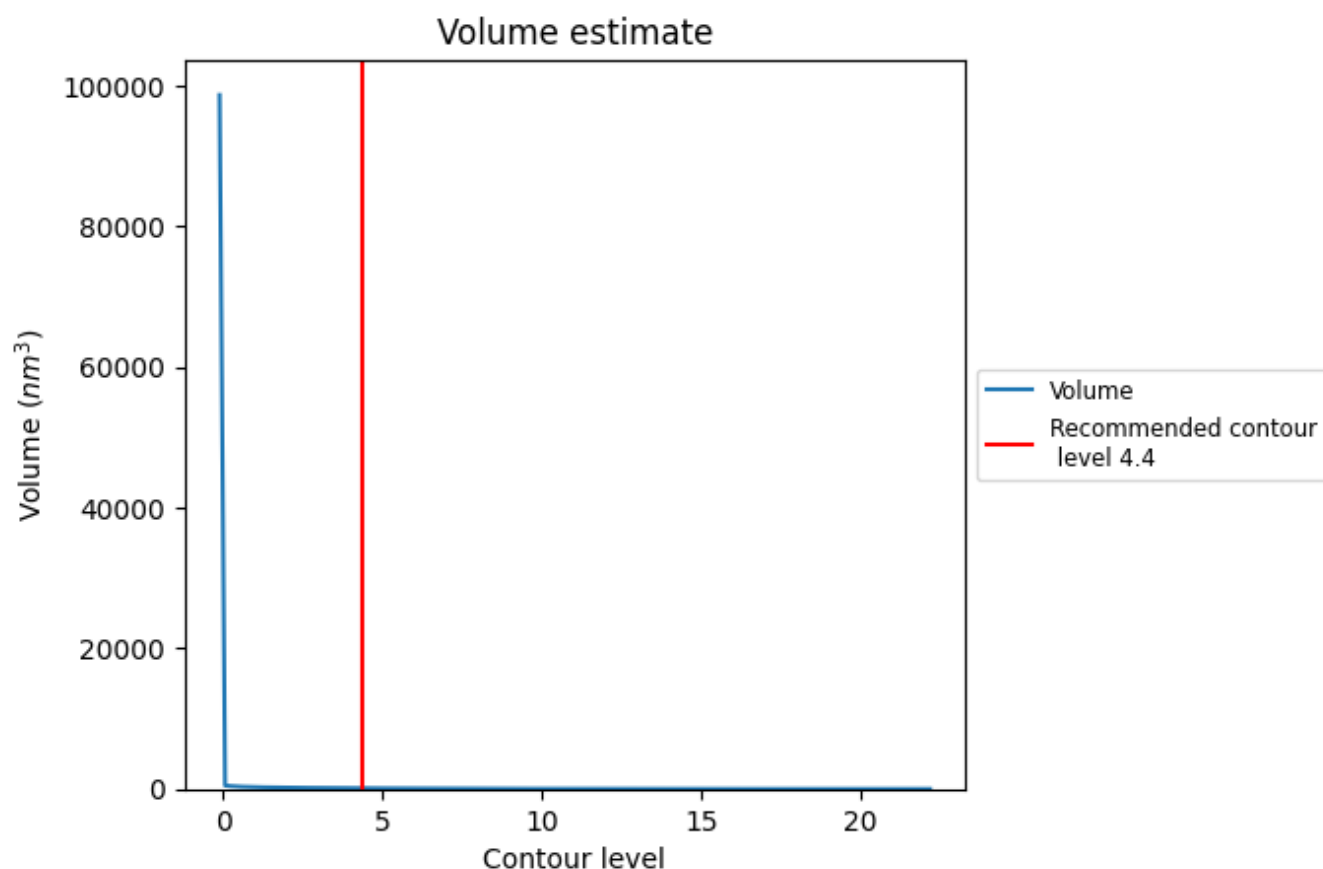
This section contains the results of statistical analysis of the map.

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.

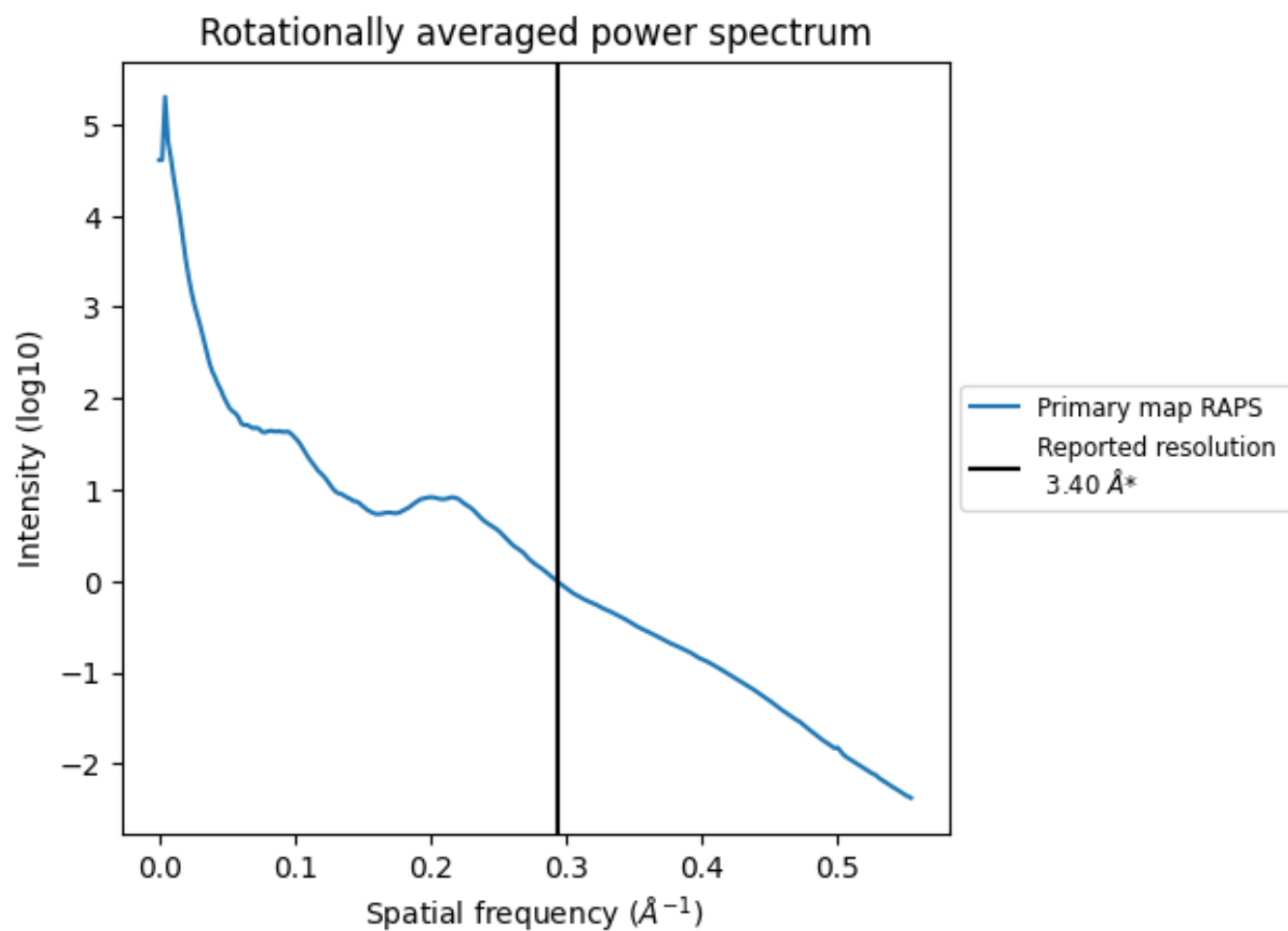
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 100 nm^3 ; this corresponds to an approximate mass of 90 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 0.294 Å⁻¹

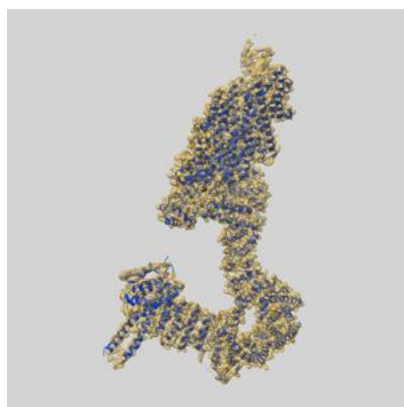
8 Fourier-Shell correlation

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

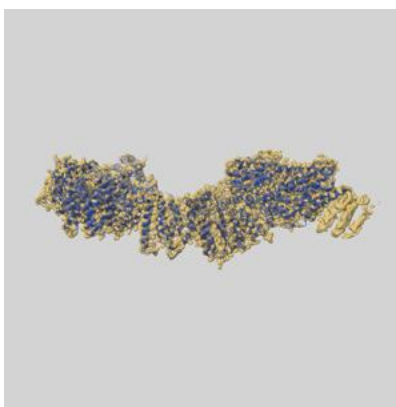
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-52297 and PDB model 9HMS. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

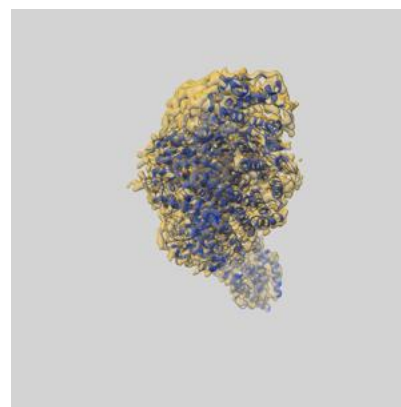
9.1 Map-model overlay [i](#)



X



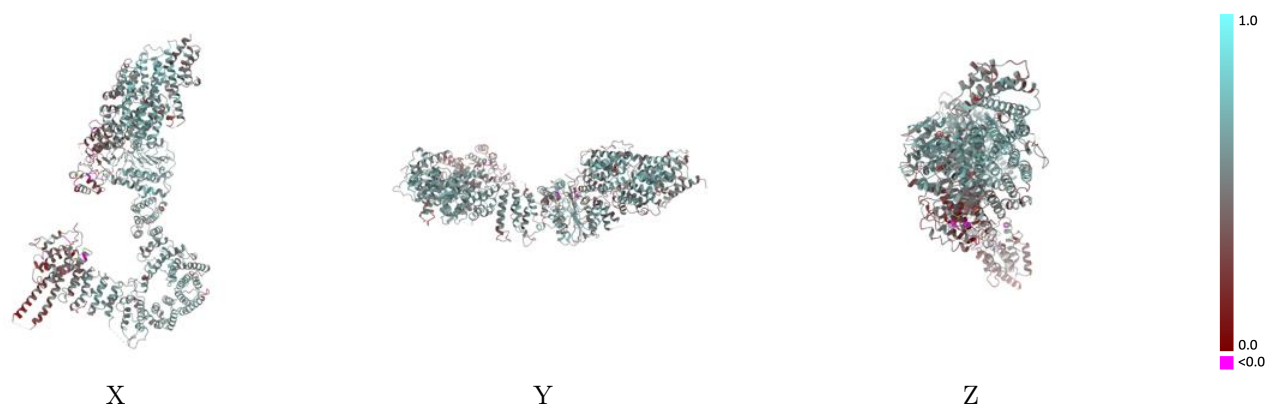
Y



Z

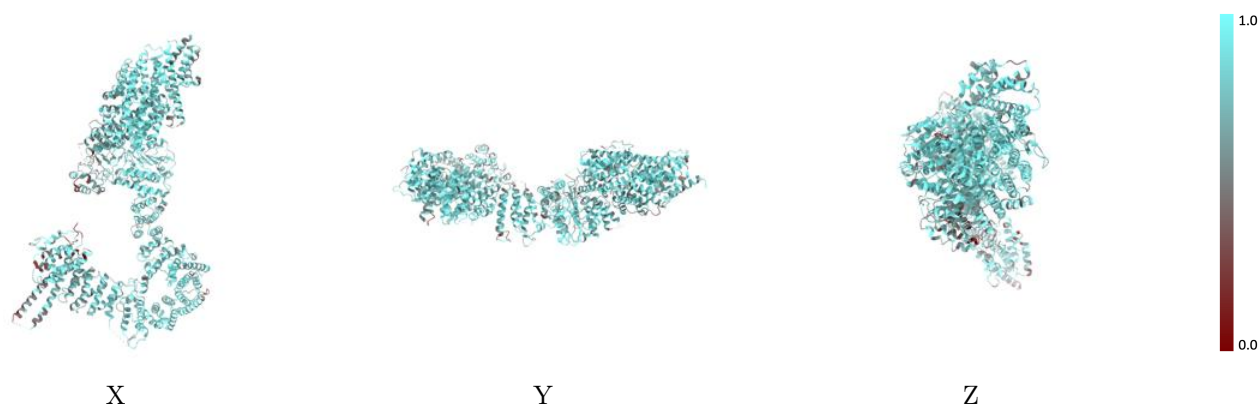
The images above show the 3D surface view of the map at the recommended contour level 4.4 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



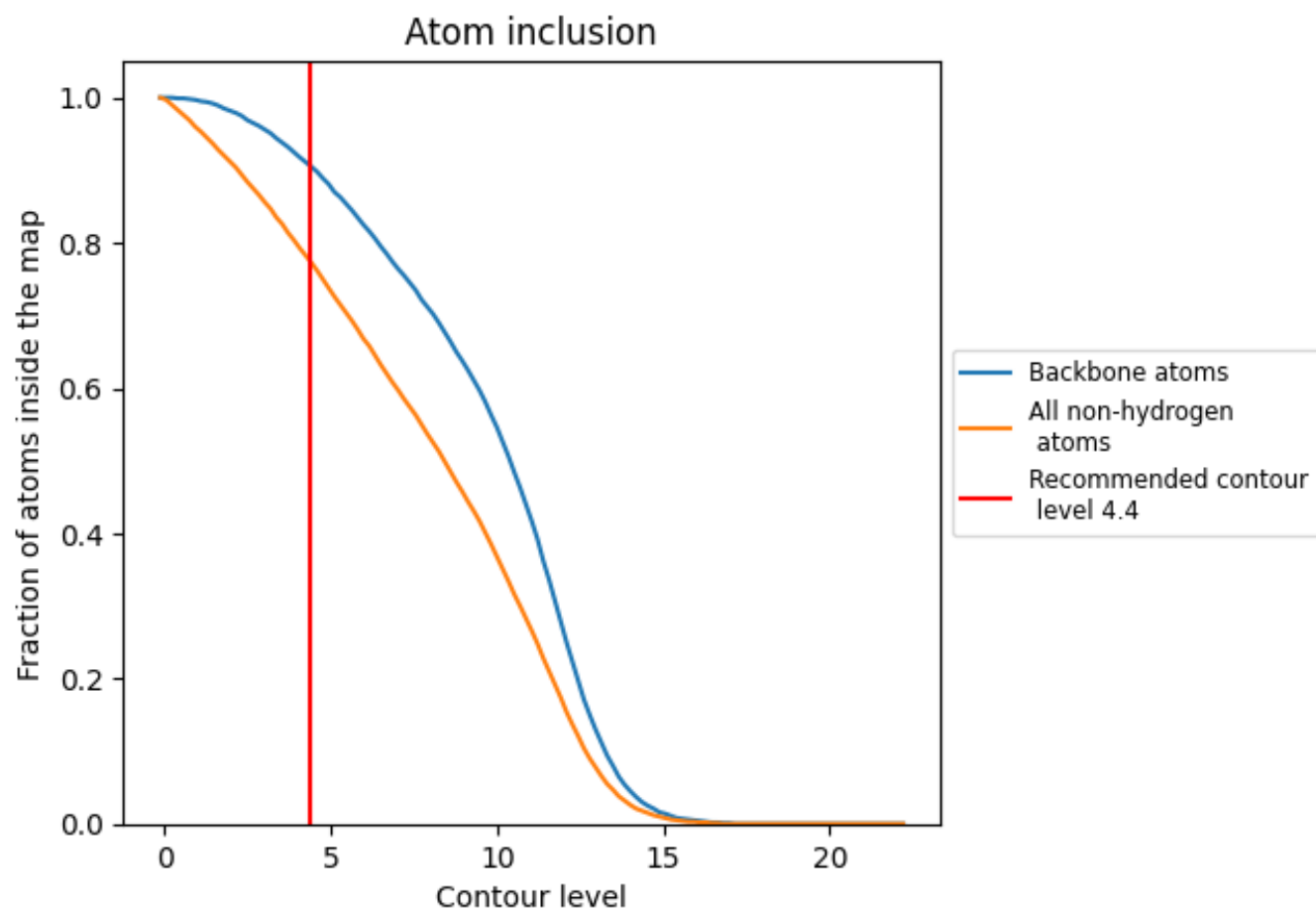
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (4.4).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 77% 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 (4.4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.7740	<div><div></div></div> 0.5010
A	<div><div></div></div> 0.7630	<div><div></div></div> 0.4940
B	<div><div></div></div> 0.6720	<div><div></div></div> 0.4640
C	<div><div></div></div> 0.7930	<div><div></div></div> 0.5090

