



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 04:56 AM UTC

PDB ID : 7LSX / pdb_00007lsx
EMDB ID : EMD-23508
Title : Cryo-EM structure of 13S proteasome core particle assembly intermediate purified from Pre3-1 proteasome mutant (G34D)
Authors : Schnell, H.M.; Walsh Jr, R.M.; Rawson, S.; Hanna, J.W.
Deposited on : 2021-02-18
Resolution : 3.61 Å(reported)
Based on initial model : 7LS6

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:

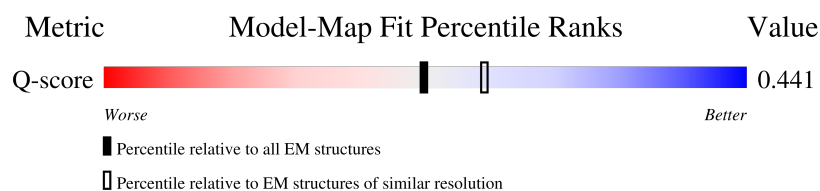
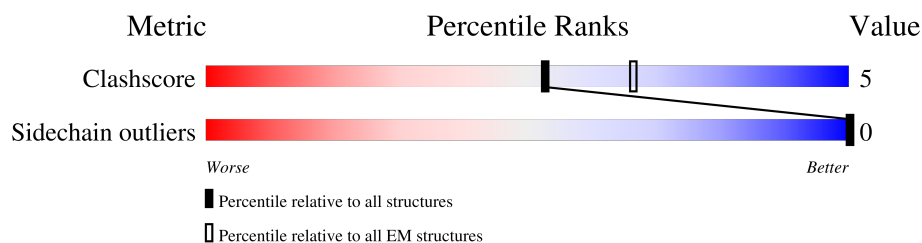
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.61 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	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11801 (3.11 - 4.10)

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	252	8% 87% 10% .
2	B	250	. 86% 13% .
3	C	258	5% 79% 12% 9%
4	D	254	8% 86% 6% 7%
5	E	260	25% 80% 16% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
6	F	234	
7	G	288	
8	H	148	
9	I	261	
10	J	205	
11	K	198	
12	O	276	
13	P	267	

2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 44393 atoms, of which 22176 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 Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	245	Total	C	H	N	O	S	0	0
			3850	1226	1922	324	370	8		

- Molecule 2 is a protein called Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	247	Total	C	H	N	O	S	0	0
			3791	1204	1900	312	372	3		

- Molecule 3 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	C	234	Total	C	H	N	O	S	0	0
			3676	1165	1840	304	364	3		

- Molecule 4 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	D	235	Total	C	H	N	O	S	0	0
			3700	1153	1859	322	362	4		

- Molecule 5 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	E	249	Total	C	H	N	O	S	0	0
			3827	1204	1903	324	388	8		

- Molecule 6 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	F	234	Total	C	H	N	O	S	0	0
			3614	1134	1811	313	351	5		

- Molecule 7 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms						AltConf	Trace
7	G	245	Total	C	H	N	O	S	0	0
			3796	1208	1895	330	359	4		

- Molecule 8 is a protein called Proteasome maturation factor UMP1.

Mol	Chain	Residues	Atoms						AltConf	Trace
8	H	91	Total	C	H	N	O	S	0	0
			1490	467	733	137	148	5		

- Molecule 9 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	I	207	Total	C	H	N	O	S	0	0
			3131	990	1563	271	302	5		

- Molecule 10 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms						AltConf	Trace
10	J	176	Total	C	H	N	O	S	0	0
			2740	890	1367	216	260	7		

- Molecule 11 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms						AltConf	Trace
11	K	182	Total	C	H	N	O	S	0	0
			2937	935	1470	247	279	6		

- Molecule 12 is a protein called Proteasome chaperone 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
12	O	240	Total	C	H	N	O	S	0	0
			3768	1216	1898	290	351	13		

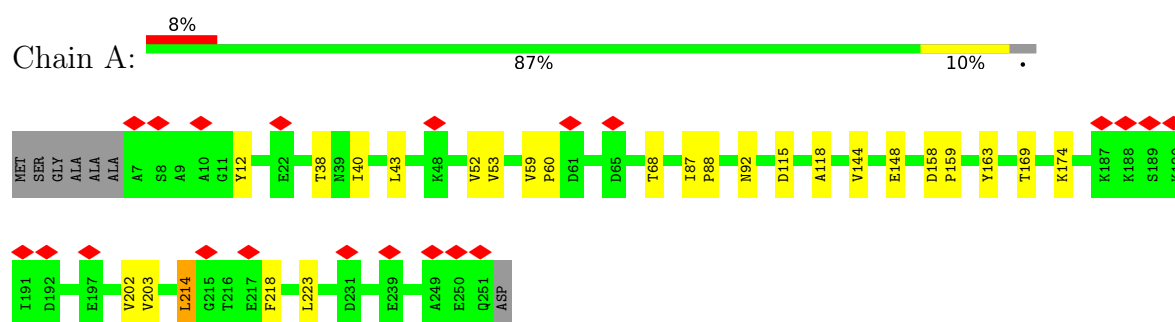
- Molecule 13 is a protein called Proteasome assembly chaperone 2.

Mol	Chain	Residues	Atoms						AltConf	Trace
13	P	253	Total	C	H	N	O	S	0	0
			4073	1333	2015	327	390	8		

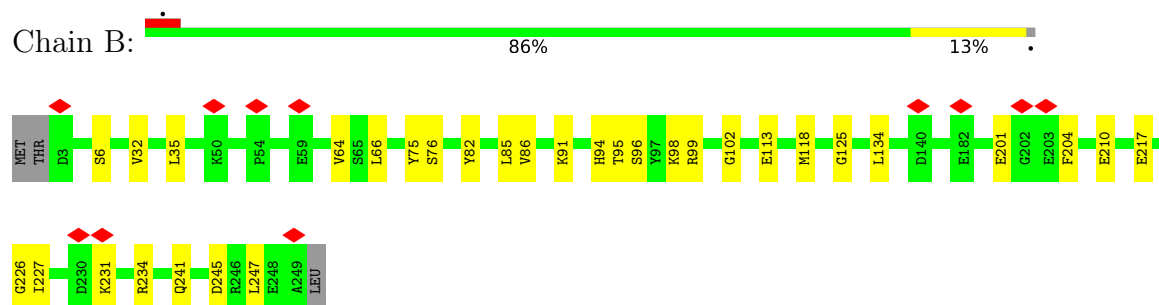
3 Residue-property plots [i](#)

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.

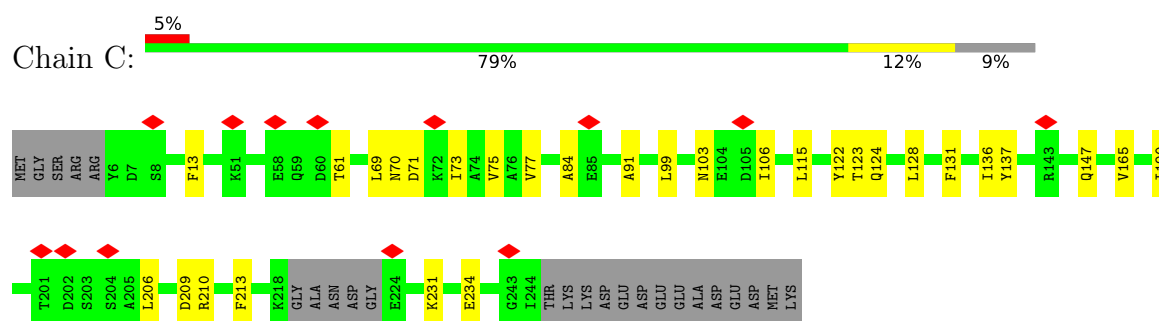
- Molecule 1: Proteasome subunit alpha type-1



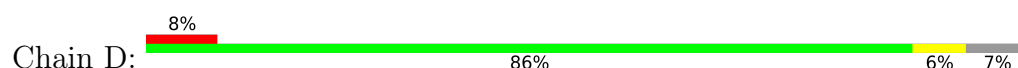
- Molecule 2: Proteasome subunit alpha type-2

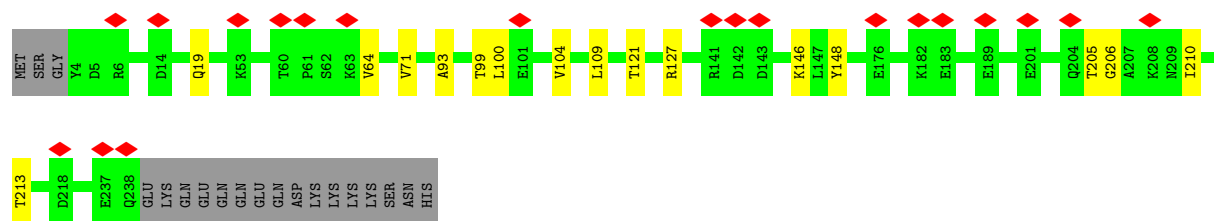


- Molecule 3: Proteasome subunit alpha type-3

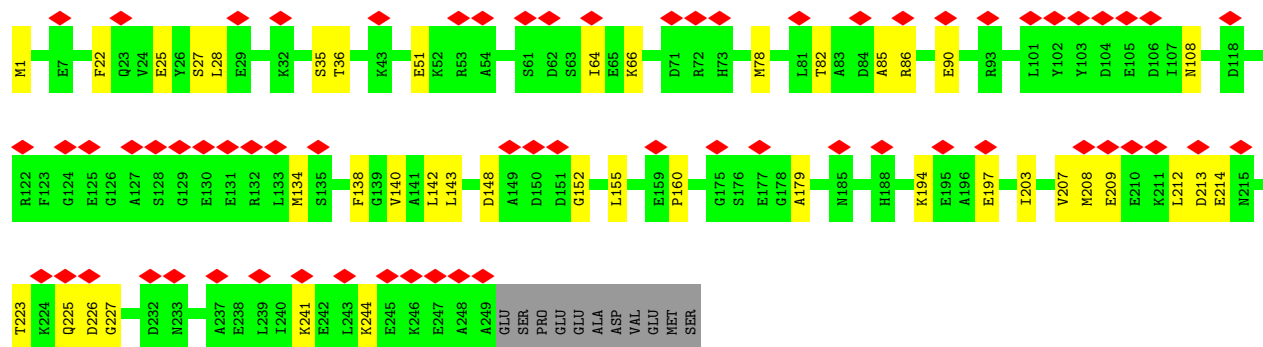
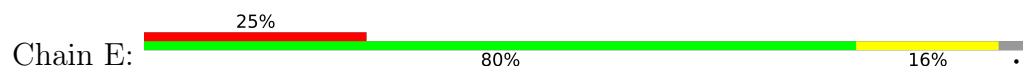


- Molecule 4: Proteasome subunit alpha type-4

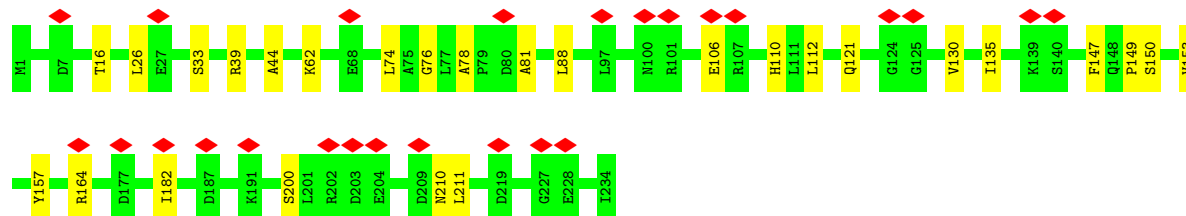
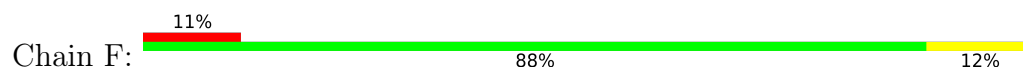




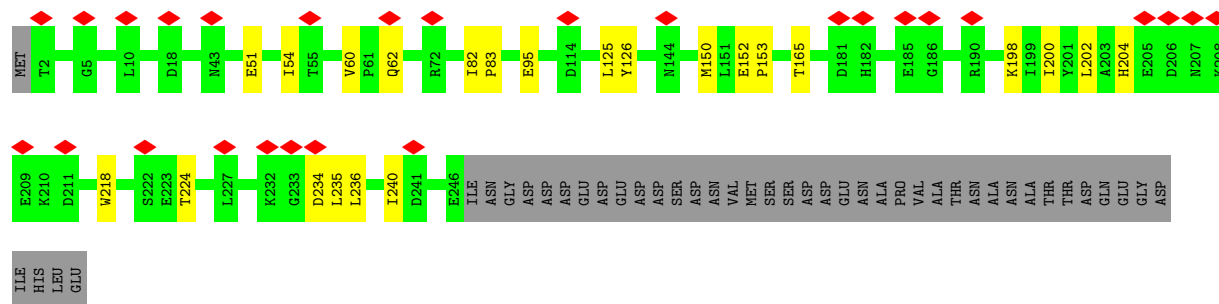
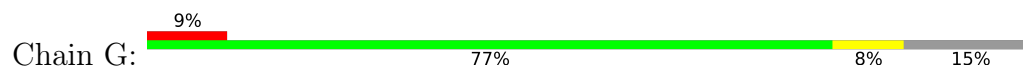
• Molecule 5: Proteasome subunit alpha type-5



• Molecule 6: Proteasome subunit alpha type-6

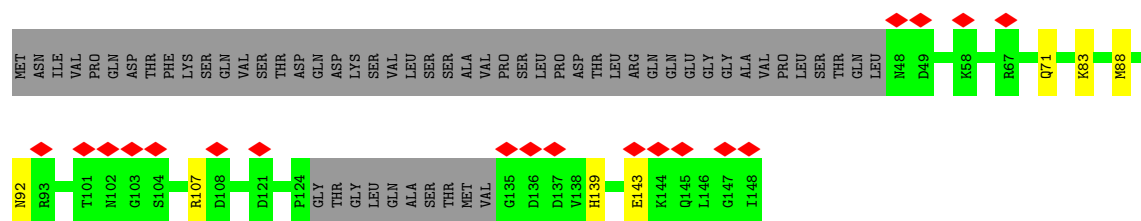


• Molecule 7: Proteasome subunit alpha type-7

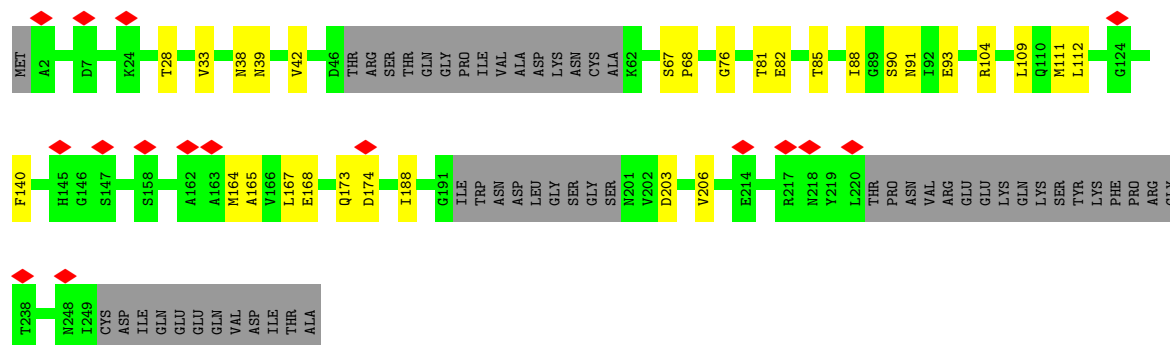


• Molecule 8: Proteasome maturation factor UMP1

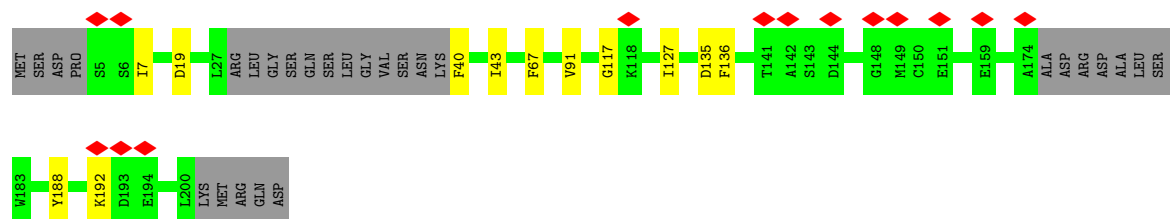
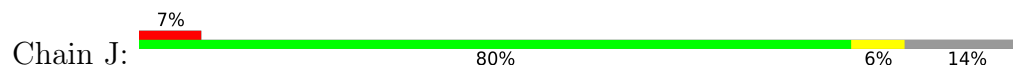




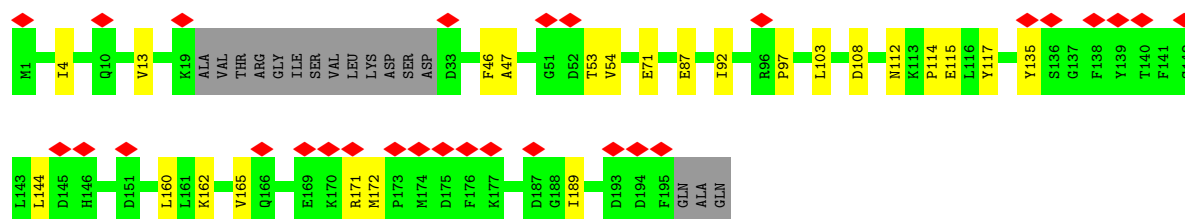
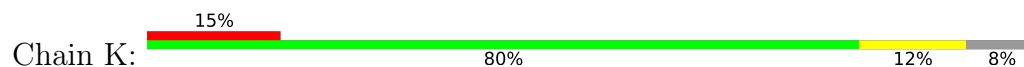
• Molecule 9: Proteasome subunit beta type-2



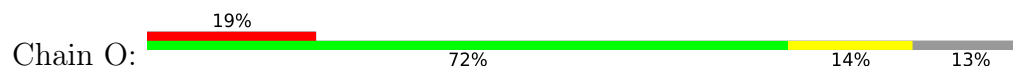
• Molecule 10: Proteasome subunit beta type-3

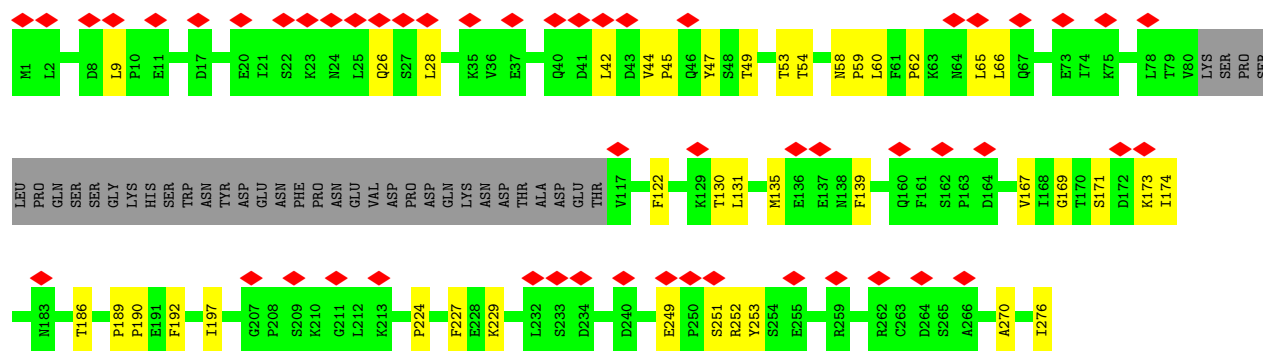


• Molecule 11: Proteasome subunit beta type-4

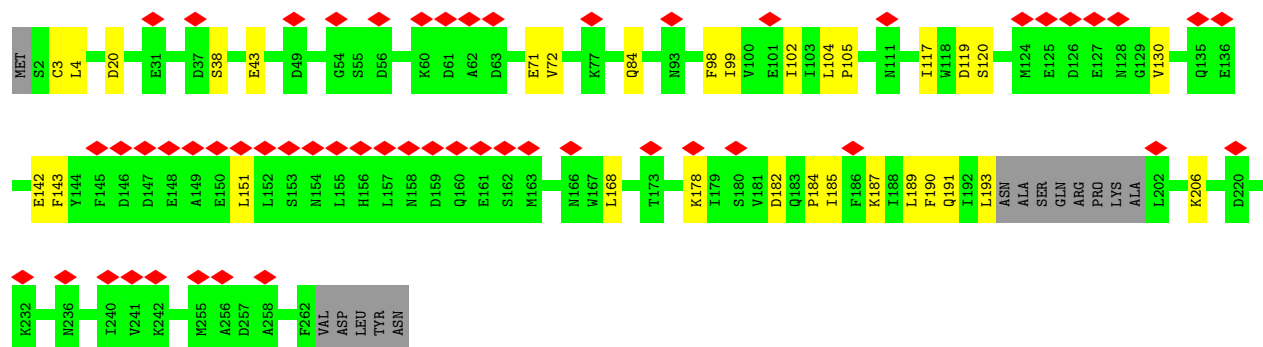
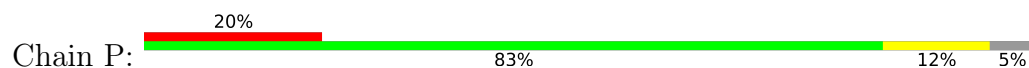


• Molecule 12: Proteasome chaperone 1





• Molecule 13: Proteasome assembly chaperone 2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	76731	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 ($e^-/\text{\AA}^2$)	55.94	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	47169	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.001	Depositor
Minimum map value	-1.005	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.051	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	381.59998, 381.59998, 381.59998	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.18	0/1966	0.38	1/2663 (0.0%)
2	B	0.14	0/1928	0.32	0/2611
3	C	0.14	0/1865	0.30	0/2525
4	D	0.13	0/1870	0.29	0/2534
5	E	0.12	0/1951	0.31	0/2628
6	F	0.13	0/1831	0.30	0/2473
7	G	0.14	0/1941	0.31	0/2621
8	H	0.12	0/770	0.31	0/1036
9	I	0.15	0/1594	0.33	0/2159
10	J	0.16	0/1400	0.32	0/1891
11	K	0.14	0/1494	0.35	0/2012
12	O	0.14	0/1914	0.33	0/2597
13	P	0.14	0/2108	0.32	0/2864
All	All	0.14	0/22632	0.32	1/30614 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	LEU	N-CA-C	-7.57	103.03	111.28

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	1928	1922	1921	18	0
2	B	1891	1900	1899	23	0
3	C	1836	1840	1838	27	0
4	D	1841	1859	1856	10	0
5	E	1924	1903	1902	32	0
6	F	1803	1811	1809	21	0
7	G	1901	1895	1892	13	0
8	H	757	733	731	6	0
9	I	1568	1563	1556	22	0
10	J	1373	1367	1364	7	0
11	K	1467	1470	1469	17	0
12	O	1870	1898	1897	30	0
13	P	2058	2015	2013	22	0
All	All	22217	22176	22147	235	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 5.

All (235) 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)
5:E:35:SER:OG	5:E:66:LYS:NZ	1.99	0.94
3:C:69:LEU:HD11	3:C:75:VAL:HG13	1.54	0.90
1:A:115:ASP:OD1	1:A:163:TYR:OH	1.94	0.85
1:A:92:ASN:OD1	8:H:83:LYS:NZ	2.15	0.79
13:P:20:ASP:OD1	13:P:84:GLN:NE2	2.19	0.76
13:P:189:LEU:HD21	13:P:193:LEU:HD12	1.68	0.76
13:P:142:GLU:OE1	13:P:206:LYS:NZ	2.19	0.75
2:B:94:HIS:ND1	9:I:90:SER:OG	2.19	0.75
7:G:51:GLU:OE2	7:G:204:HIS:ND1	2.23	0.71
9:I:33:VAL:HG22	9:I:188:ILE:HD11	1.72	0.71
1:A:174:LYS:HD3	1:A:214:LEU:HD13	1.71	0.71
12:O:249:GLU:CD	12:O:251:SER:HG	1.99	0.70
13:P:71:GLU:N	13:P:71:GLU:OE1	2.24	0.70
3:C:147:GLN:N	3:C:147:GLN:OE1	2.28	0.67
11:K:135:TYR:CE2	11:K:172:MET:HE3	2.31	0.66
3:C:69:LEU:CD1	3:C:75:VAL:HG13	2.24	0.66
5:E:28:LEU:HD12	12:O:270:ALA:HB2	1.78	0.66
12:O:49:THR:CG2	12:O:130:THR:HG22	2.26	0.65
1:A:53:VAL:CG1	1:A:144:VAL:HG11	2.28	0.64
5:E:22:PHE:O	5:E:25:GLU:N	2.31	0.64
6:F:106:GLU:OE2	6:F:110:HIS:ND1	2.31	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:135:TYR:CZ	11:K:172:MET:HE3	2.34	0.63
5:E:64:ILE:HG23	5:E:64:ILE:O	1.99	0.62
2:B:85:LEU:CD2	2:B:118:MET:HE3	2.30	0.62
13:P:119:ASP:OD1	13:P:120:SER:N	2.34	0.61
3:C:136:ILE:HD11	3:C:165:VAL:CG2	2.31	0.60
11:K:115:GLU:OE1	11:K:117:TYR:OH	2.19	0.60
3:C:91:ALA:HB2	3:C:115:LEU:HD11	1.84	0.60
3:C:122:TYR:CD1	3:C:128:LEU:HD13	2.37	0.60
7:G:54:ILE:HG23	7:G:54:ILE:O	2.01	0.60
5:E:85:ALA:HB2	5:E:140:VAL:HG11	1.83	0.59
7:G:234:ASP:OD1	7:G:235:LEU:N	2.36	0.59
6:F:33:SER:OG	6:F:62:LYS:NZ	2.28	0.58
6:F:78:ALA:HA	12:O:276:ILE:HD11	1.85	0.58
6:F:44:ALA:C	6:F:135:ILE:HD11	2.29	0.58
1:A:218:PHE:HE1	1:A:223:LEU:HD13	1.70	0.57
6:F:182:ILE:HG22	6:F:182:ILE:O	2.04	0.57
2:B:204:PHE:CE2	2:B:247:LEU:HD11	2.40	0.57
10:J:135:ASP:OD1	10:J:136:PHE:N	2.37	0.56
9:I:164:MET:O	9:I:168:GLU:OE1	2.23	0.56
3:C:13:PHE:N	4:D:19:GLN:OE1	2.39	0.56
13:P:189:LEU:CD2	13:P:193:LEU:HD12	2.33	0.56
3:C:136:ILE:HD11	3:C:165:VAL:HG21	1.89	0.54
3:C:69:LEU:HD11	3:C:75:VAL:CG1	2.31	0.54
9:I:38:ASN:OD1	9:I:39:ASN:N	2.40	0.54
6:F:210:ASN:OD1	6:F:211:LEU:N	2.39	0.54
12:O:186:THR:O	12:O:186:THR:HG22	2.07	0.54
11:K:171:ARG:O	11:K:172:MET:HE2	2.08	0.54
1:A:218:PHE:CE1	1:A:223:LEU:HD13	2.42	0.54
9:I:28:THR:OG1	9:I:76:GLY:N	2.41	0.54
6:F:74:LEU:HD22	6:F:81:ALA:HB1	1.89	0.53
11:K:4:ILE:HG22	11:K:103:LEU:HD12	1.90	0.53
11:K:4:ILE:CD1	11:K:47:ALA:HB2	2.37	0.53
5:E:223:THR:O	5:E:227:GLY:N	2.36	0.53
5:E:108:ASN:ND2	5:E:148:ASP:OD2	2.39	0.53
1:A:68:THR:HG22	1:A:68:THR:O	2.09	0.52
12:O:28:LEU:HD11	13:P:189:LEU:HD13	1.89	0.52
2:B:64:VAL:HG22	2:B:210:GLU:HG2	1.91	0.52
3:C:124:GLN:OE1	4:D:127:ARG:N	2.43	0.52
12:O:49:THR:HG23	12:O:130:THR:HG22	1.90	0.52
5:E:28:LEU:HD12	12:O:270:ALA:CB	2.39	0.52
1:A:87:ILE:N	1:A:88:PRO:HD2	2.24	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:117:GLY:O	10:J:192:LYS:NZ	2.43	0.52
3:C:231:LYS:N	3:C:234:GLU:OE2	2.36	0.52
5:E:27:SER:OG	5:E:138:PHE:O	2.29	0.51
11:K:189:ILE:HD12	11:K:189:ILE:H	1.76	0.51
5:E:179:ALA:HA	5:E:203:ILE:HD11	1.93	0.51
10:J:7:ILE:HD11	10:J:40:PHE:CD2	2.46	0.51
2:B:91:LYS:O	2:B:95:THR:HG22	2.10	0.51
7:G:150:MET:HE3	7:G:165:THR:HG22	1.93	0.50
12:O:276:ILE:HD12	12:O:276:ILE:H	1.76	0.50
13:P:187:LYS:O	13:P:190:PHE:N	2.44	0.50
3:C:122:TYR:CE1	3:C:128:LEU:HD13	2.46	0.50
2:B:113:GLU:OE1	2:B:113:GLU:HA	2.12	0.50
11:K:4:ILE:CG2	11:K:103:LEU:HD12	2.41	0.50
2:B:98:LYS:O	2:B:102:GLY:N	2.39	0.50
1:A:52:VAL:HG21	1:A:203:VAL:HG22	1.92	0.50
2:B:82:TYR:O	2:B:86:VAL:HG23	2.12	0.50
12:O:122:PHE:CE2	12:O:135:MET:HE1	2.47	0.50
4:D:71:VAL:HG11	4:D:109:LEU:CD2	2.43	0.49
6:F:164:ARG:NH1	6:F:200:SER:O	2.42	0.49
5:E:28:LEU:HD21	5:E:160:PRO:HD2	1.94	0.49
5:E:241:LYS:HA	5:E:244:LYS:HG2	1.94	0.49
6:F:88:LEU:HD12	6:F:112:LEU:HD12	1.94	0.49
6:F:39:ARG:N	6:F:157:TYR:O	2.46	0.49
13:P:182:ASP:O	13:P:185:ILE:HG22	2.13	0.49
5:E:78:MET:HE1	5:E:82:THR:HG22	1.95	0.48
9:I:93:GLU:OE1	9:I:93:GLU:HA	2.12	0.48
1:A:118:ALA:HB3	1:A:163:TYR:HE1	1.79	0.48
2:B:118:MET:HE1	2:B:134:LEU:HD13	1.95	0.48
11:K:108:ASP:O	11:K:112:ASN:N	2.47	0.48
6:F:121:GLN:OE1	6:F:121:GLN:HA	2.13	0.48
9:I:173:GLN:O	9:I:174:ASP:OD1	2.32	0.48
9:I:88:ILE:HD12	9:I:112:LEU:HD23	1.96	0.48
8:H:71:GLN:OE1	8:H:71:GLN:HA	2.13	0.47
10:J:19:ASP:OD1	10:J:192:LYS:HD3	2.14	0.47
6:F:76:GLY:O	12:O:276:ILE:HD12	2.14	0.47
13:P:72:VAL:HG23	13:P:72:VAL:O	2.15	0.47
5:E:86:ARG:NH1	5:E:90:GLU:OE2	2.46	0.47
11:K:144:LEU:HD23	11:K:160:LEU:HD21	1.96	0.47
2:B:241:GLN:NE2	2:B:245:ASP:OD2	2.43	0.47
7:G:125:LEU:HD23	7:G:126:TYR:CZ	2.49	0.47
13:P:182:ASP:OD1	13:P:182:ASP:N	2.47	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:99:ARG:HA	2:B:99:ARG:NE	2.30	0.47
3:C:136:ILE:HD11	3:C:165:VAL:HG23	1.96	0.47
7:G:95:GLU:OE1	7:G:95:GLU:HA	2.14	0.47
2:B:66:LEU:H	2:B:66:LEU:HD23	1.79	0.47
12:O:65:LEU:HD12	12:O:252:ARG:NH1	2.30	0.47
12:O:227:PHE:O	12:O:229:LYS:NZ	2.41	0.47
9:I:174:ASP:OD1	9:I:174:ASP:C	2.58	0.47
9:I:81:THR:O	9:I:85:THR:OG1	2.29	0.46
12:O:26:GLN:O	12:O:26:GLN:NE2	2.43	0.46
7:G:60:VAL:HG22	7:G:62:GLN:OE1	2.15	0.46
9:I:164:MET:O	9:I:167:LEU:N	2.48	0.46
10:J:67:PHE:CZ	10:J:91:VAL:HG23	2.51	0.46
12:O:139:PHE:CE1	12:O:224:PRO:HB2	2.50	0.46
12:O:58:ASN:N	12:O:59:PRO:CD	2.79	0.46
1:A:158:ASP:HB2	1:A:159:PRO:HD2	1.98	0.46
5:E:142:LEU:C	5:E:142:LEU:HD12	2.40	0.46
9:I:109:LEU:C	9:I:109:LEU:HD23	2.41	0.46
2:B:217:GLU:OE2	2:B:234:ARG:NH2	2.49	0.46
9:I:203:ASP:OD1	9:I:203:ASP:N	2.46	0.46
12:O:42:LEU:HD22	12:O:44:VAL:HG22	1.98	0.46
3:C:61:THR:HG22	3:C:61:THR:O	2.16	0.45
4:D:93:ALA:HA	4:D:104:VAL:HG11	1.98	0.45
6:F:16:THR:O	6:F:16:THR:HG22	2.15	0.45
5:E:134:MET:HE1	5:E:138:PHE:CE1	2.52	0.45
5:E:148:ASP:O	5:E:152:GLY:N	2.43	0.45
13:P:38:SER:OG	13:P:43:GLU:OE2	2.28	0.45
1:A:53:VAL:HG13	1:A:144:VAL:HG11	1.97	0.45
7:G:236:LEU:O	7:G:240:ILE:HG12	2.17	0.45
4:D:205:THR:HG22	4:D:206:GLY:N	2.32	0.45
3:C:69:LEU:HD13	3:C:73:ILE:HG22	1.98	0.45
6:F:26:LEU:HD22	6:F:149:PRO:HG2	1.97	0.45
7:G:198:LYS:O	7:G:202:LEU:HD23	2.17	0.45
12:O:62:PRO:O	12:O:66:LEU:HD23	2.16	0.45
12:O:171:SER:HB3	12:O:174:ILE:HD12	1.98	0.45
13:P:104:LEU:HB2	13:P:105:PRO:HD3	1.99	0.45
12:O:60:LEU:O	12:O:253:TYR:OH	2.34	0.44
13:P:3:CYS:SG	13:P:4:LEU:N	2.90	0.44
2:B:95:THR:HG23	2:B:96:SER:N	2.32	0.44
3:C:122:TYR:CD2	3:C:131:PHE:CE2	3.06	0.44
2:B:35:LEU:C	2:B:35:LEU:HD12	2.43	0.44
11:K:13:VAL:HG23	11:K:114:PRO:HB2	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:O:47:TYR:CE2	12:O:131:LEU:HD21	2.53	0.44
13:P:117:ILE:HG21	13:P:187:LYS:HE2	2.00	0.44
3:C:75:VAL:HG12	3:C:137:TYR:HA	1.99	0.44
3:C:91:ALA:CB	3:C:115:LEU:HD11	2.47	0.44
6:F:44:ALA:HB1	6:F:135:ILE:HD11	1.99	0.44
11:K:71:GLU:HA	11:K:71:GLU:OE1	2.18	0.44
4:D:121:THR:HG22	4:D:121:THR:O	2.18	0.44
5:E:143:LEU:HD12	5:E:155:LEU:HD21	2.00	0.44
9:I:42:VAL:HG12	9:I:206:VAL:HG23	2.00	0.44
12:O:167:VAL:O	12:O:167:VAL:HG13	2.17	0.44
4:D:99:THR:C	4:D:100:LEU:HD12	2.43	0.44
13:P:104:LEU:HD21	13:P:191:GLN:HG3	1.99	0.44
3:C:61:THR:O	3:C:61:THR:CG2	2.66	0.43
7:G:82:ILE:N	7:G:83:PRO:HD2	2.33	0.43
8:H:88:MET:HE1	8:H:107:ARG:HG3	2.00	0.43
5:E:36:THR:HG21	5:E:207:VAL:HG11	1.99	0.43
1:A:52:VAL:HG11	1:A:202:VAL:HG12	1.99	0.43
5:E:51:GLU:OE1	5:E:208:MET:HE2	2.19	0.43
9:I:67:SER:O	9:I:68:PRO:C	2.61	0.43
1:A:43:LEU:HD12	1:A:169:THR:O	2.18	0.43
1:A:148:GLU:OE2	9:I:104:ARG:NH1	2.45	0.43
6:F:44:ALA:CB	6:F:135:ILE:HD11	2.49	0.43
8:H:88:MET:HE3	8:H:92:ASN:HD21	1.82	0.43
9:I:164:MET:O	9:I:165:ALA:C	2.59	0.43
2:B:6:SER:OG	2:B:125:GLY:O	2.30	0.43
5:E:25:GLU:OE1	12:O:270:ALA:HB3	2.18	0.43
9:I:82:GLU:OE1	9:I:82:GLU:HA	2.19	0.43
5:E:155:LEU:C	5:E:155:LEU:HD23	2.43	0.43
5:E:213:ASP:OD1	5:E:214:GLU:N	2.52	0.43
7:G:218:TRP:CZ3	7:G:224:THR:HG23	2.54	0.43
11:K:162:LYS:O	11:K:165:VAL:HG12	2.19	0.43
5:E:1:MET:HG3	5:E:1:MET:O	2.19	0.42
12:O:173:LYS:HG3	12:O:192:PHE:CZ	2.54	0.42
13:P:99:ILE:HD12	13:P:191:GLN:NE2	2.34	0.42
2:B:85:LEU:HD21	2:B:118:MET:HE3	2.01	0.42
1:A:12:TYR:CE1	12:O:9:LEU:HD21	2.55	0.42
3:C:123:THR:O	3:C:123:THR:HG22	2.19	0.42
3:C:206:LEU:O	3:C:206:LEU:HG	2.19	0.42
12:O:53:THR:HG22	12:O:54:THR:O	2.20	0.42
1:A:38:THR:CG2	1:A:40:ILE:HG22	2.49	0.42
3:C:77:VAL:HG21	3:C:84:ALA:CB	2.50	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:91:ASN:CG	9:I:111:MET:HE1	2.45	0.42
5:E:64:ILE:O	5:E:64:ILE:CG2	2.67	0.42
5:E:209:GLU:OE1	5:E:209:GLU:N	2.49	0.42
6:F:147:PHE:CD2	6:F:153:VAL:HG22	2.54	0.42
10:J:127:ILE:O	10:J:127:ILE:HG22	2.20	0.42
1:A:59:VAL:HB	1:A:60:PRO:HA	2.02	0.42
5:E:225:GLN:HG2	5:E:226:ASP:N	2.35	0.41
7:G:51:GLU:N	7:G:200:ILE:HD11	2.35	0.41
11:K:92:ILE:HD11	11:K:97:PRO:HB3	2.02	0.41
5:E:134:MET:HE1	5:E:138:PHE:CZ	2.55	0.41
6:F:26:LEU:HD21	6:F:150:SER:HB2	2.01	0.41
13:P:130:VAL:HG13	13:P:178:LYS:O	2.19	0.41
4:D:146:LYS:NZ	4:D:148:TYR:OH	2.43	0.41
9:I:88:ILE:HD12	9:I:112:LEU:CD2	2.50	0.41
6:F:164:ARG:O	6:F:200:SER:OG	2.38	0.41
11:K:53:THR:HG23	11:K:54:VAL:N	2.36	0.41
2:B:64:VAL:HG22	2:B:210:GLU:CG	2.50	0.41
3:C:106:ILE:O	3:C:106:ILE:HG23	2.20	0.41
9:I:109:LEU:HD12	9:I:140:PHE:CB	2.50	0.41
11:K:4:ILE:HD11	11:K:46:PHE:O	2.20	0.41
12:O:169:GLY:HA3	12:O:197:ILE:HG21	2.03	0.41
3:C:70:ASN:OD1	3:C:71:ASP:N	2.52	0.41
5:E:208:MET:HE1	5:E:212:LEU:HD13	2.02	0.41
2:B:201:GLU:N	2:B:201:GLU:OE1	2.53	0.41
2:B:32:VAL:O	2:B:76:SER:OG	2.35	0.41
5:E:143:LEU:HD23	5:E:143:LEU:H	1.85	0.41
6:F:130:VAL:HG13	6:F:130:VAL:O	2.21	0.41
7:G:152:GLU:HB3	7:G:153:PRO:HD2	2.02	0.41
12:O:189:PRO:N	12:O:190:PRO:CD	2.84	0.41
3:C:190:ILE:HG12	3:C:213:PHE:HZ	1.86	0.40
4:D:210:ILE:O	4:D:210:ILE:HG23	2.21	0.40
2:B:226:GLY:O	2:B:227:ILE:C	2.63	0.40
6:F:88:LEU:HD12	6:F:112:LEU:CD1	2.50	0.40
11:K:87:GLU:HA	11:K:87:GLU:OE1	2.22	0.40
2:B:75:TYR:HB3	2:B:82:TYR:CD1	2.56	0.40
3:C:99:LEU:O	3:C:103:ASN:HA	2.21	0.40
3:C:209:ASP:OD1	3:C:210:ARG:N	2.55	0.40
4:D:64:VAL:HG11	4:D:213:THR:OG1	2.21	0.40
5:E:1:MET:HE2	12:O:189:PRO:O	2.21	0.40
5:E:194:LYS:O	5:E:197:GLU:HG2	2.21	0.40
9:I:109:LEU:HD12	9:I:140:PHE:CG	2.57	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:P:151:LEU:C	13:P:151:LEU:HD23	2.46	0.40
2:B:231:LYS:N	2:B:231:LYS:HD2	2.36	0.40
12:O:44:VAL:N	12:O:45:PRO:CD	2.84	0.40
8:H:88:MET:HE3	8:H:92:ASN:ND2	2.37	0.40
8:H:139:HIS:O	8:H:143:GLU:HG2	2.21	0.40
10:J:43:ILE:HG21	10:J:188:TYR:CD1	2.57	0.40
13:P:98:PHE:CD1	13:P:102:ILE:HD12	2.57	0.40
13:P:143:PHE:HB2	13:P:168:LEU:HD23	2.02	0.40
13:P:184:PRO:HA	13:P:187:LYS:HE3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	207/210 (99%)	207 (100%)	0	100	100
2	B	206/209 (99%)	206 (100%)	0	100	100
3	C	197/216 (91%)	197 (100%)	0	100	100
4	D	208/226 (92%)	208 (100%)	0	100	100
5	E	205/215 (95%)	205 (100%)	0	100	100
6	F	193/193 (100%)	193 (100%)	0	100	100
7	G	202/239 (84%)	202 (100%)	0	100	100
8	H	86/136 (63%)	86 (100%)	0	100	100
9	I	167/214 (78%)	167 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
10	J	149/173 (86%)	149 (100%)	0	100	100
11	K	162/175 (93%)	162 (100%)	0	100	100
12	O	217/251 (86%)	217 (100%)	0	100	100
13	P	232/244 (95%)	232 (100%)	0	100	100
All	All	2431/2701 (90%)	2431 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	209	HIS
3	C	94	HIS
5	E	100	ASN
5	E	154	GLN
5	E	168	ASN
7	G	123	HIS
7	G	170	GLN
7	G	237	GLN
8	H	64	GLN
9	I	122	HIS
11	K	78	GLN
11	K	101	ASN
11	K	133	HIS
12	O	157	GLN
13	P	154	ASN
13	P	158	ASN
13	P	213	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](#)

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.

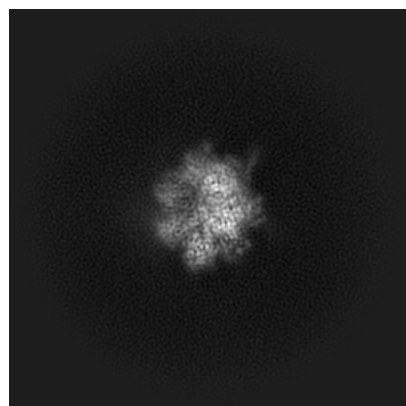
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23508. These allow visual inspection of the internal detail of the map and identification of artifacts.

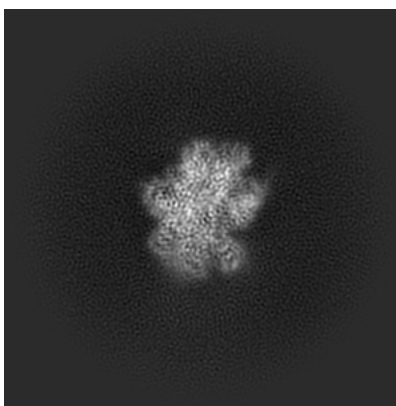
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

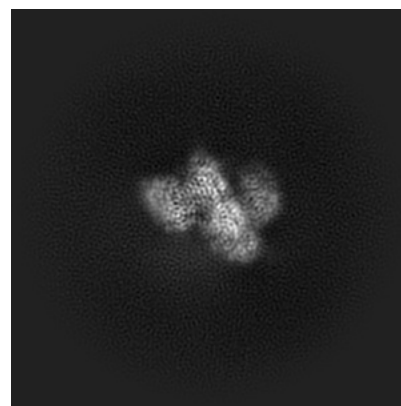
6.1.1 Primary map



X

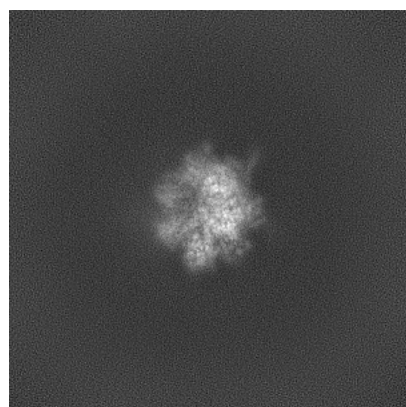


Y

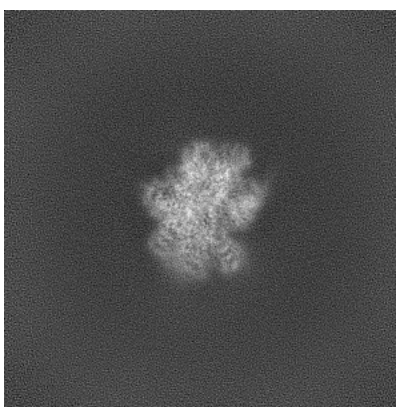


Z

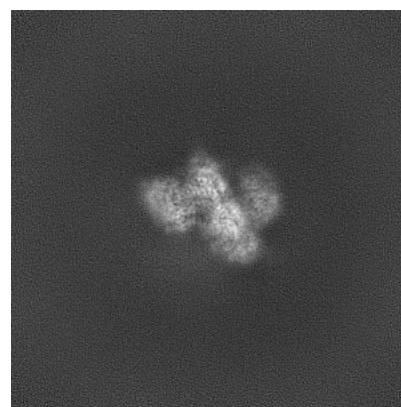
6.1.2 Raw map



X



Y

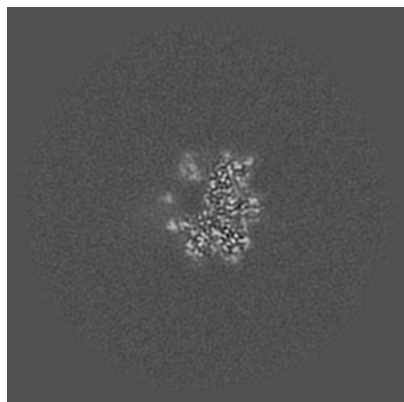


Z

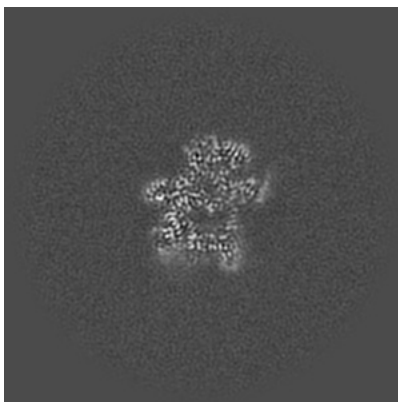
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

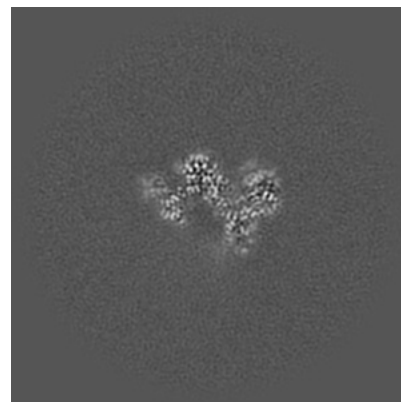
6.2.1 Primary map



X Index: 180

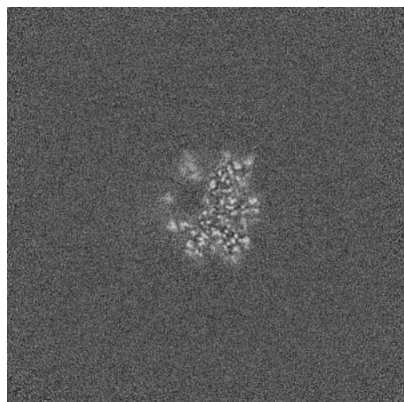


Y Index: 180

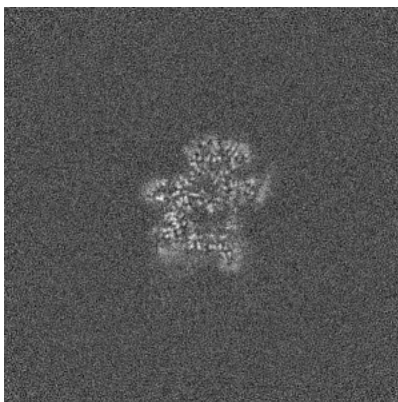


Z Index: 180

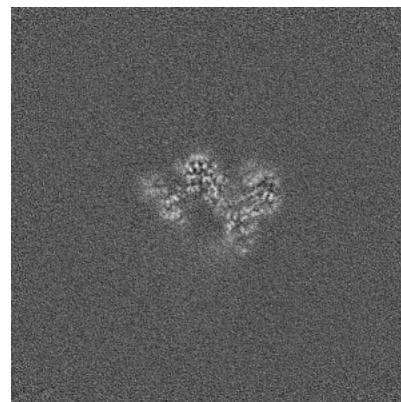
6.2.2 Raw map



X Index: 180



Y Index: 180

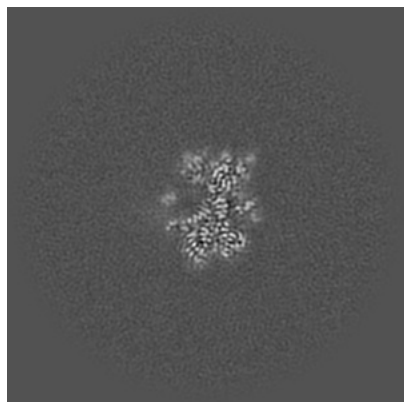


Z Index: 180

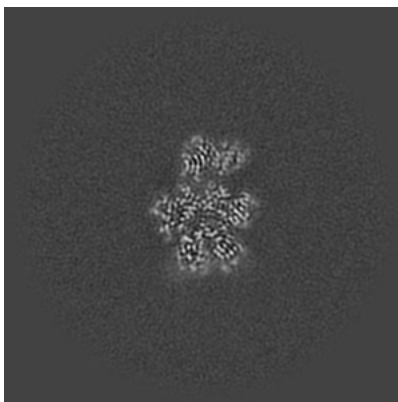
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

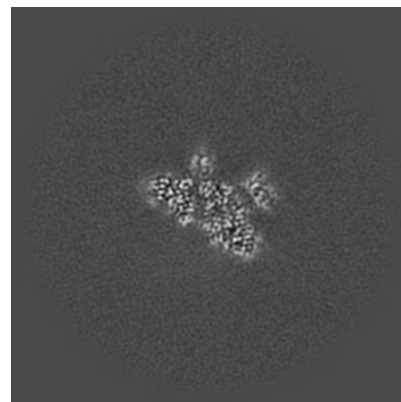
6.3.1 Primary map



X Index: 183

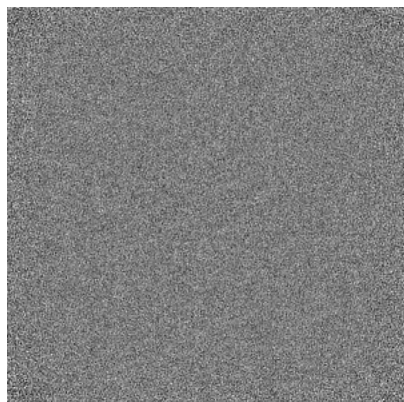


Y Index: 193

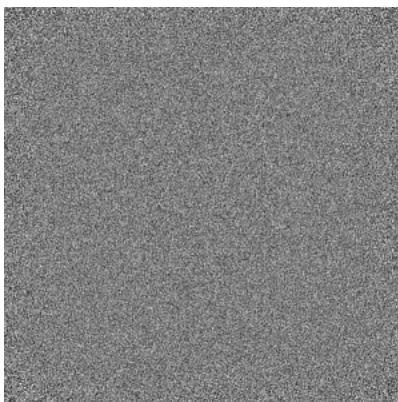


Z Index: 166

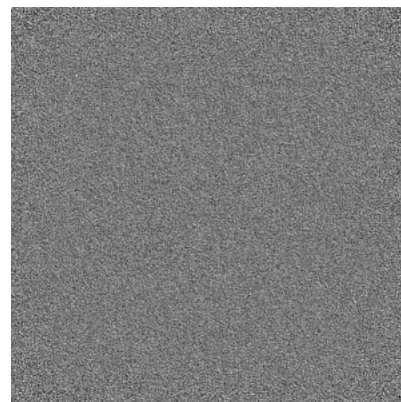
6.3.2 Raw map



X Index: 0



Y Index: 0

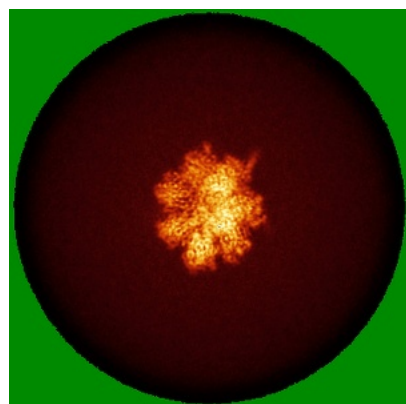


Z Index: 0

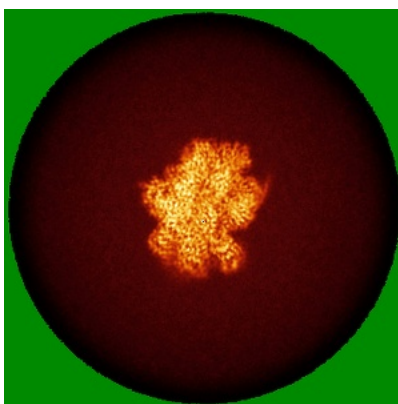
The images above show the largest variance slices of the map in three orthogonal directions.

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

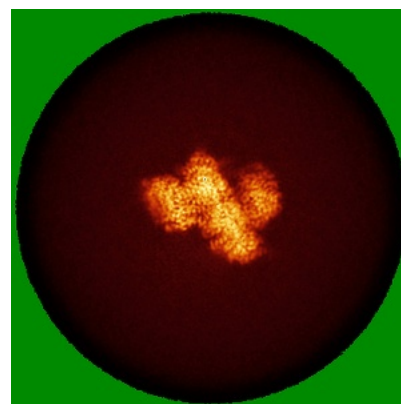
6.4.1 Primary map



X

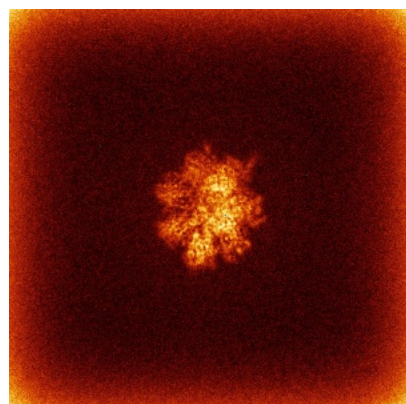


Y

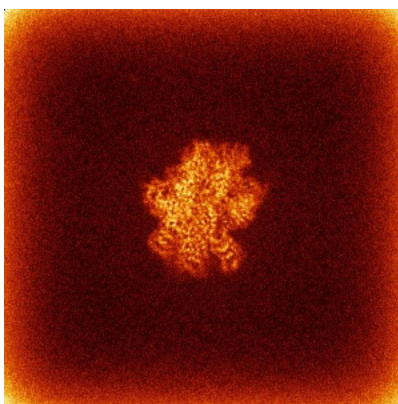


Z

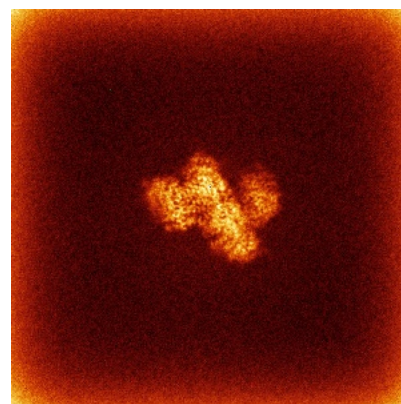
6.4.2 Raw map



X



Y

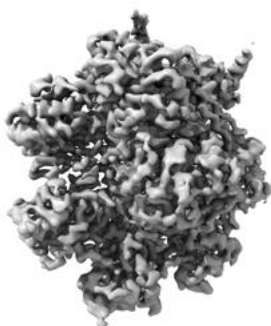


Z

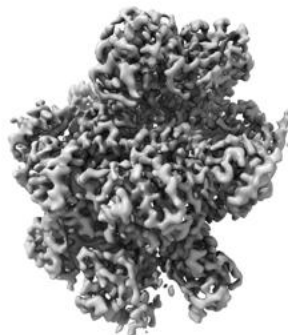
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

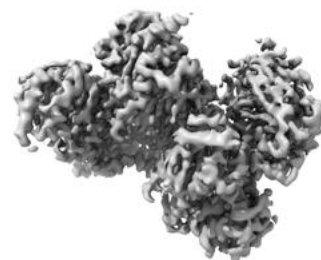
6.5.1 Primary map



X



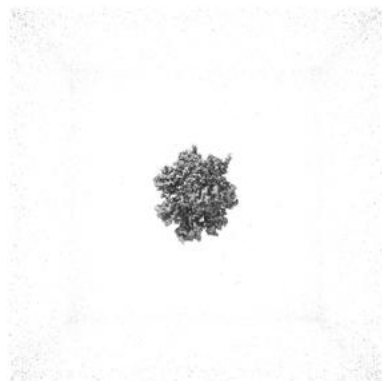
Y



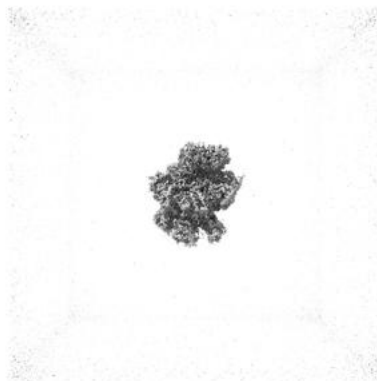
Z

The images above show the 3D surface view of the map at the recommended contour level 0.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

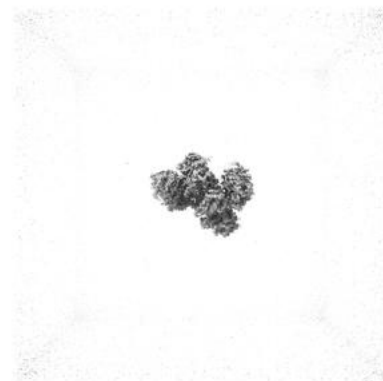
6.5.2 Raw map



X



Y



Z

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

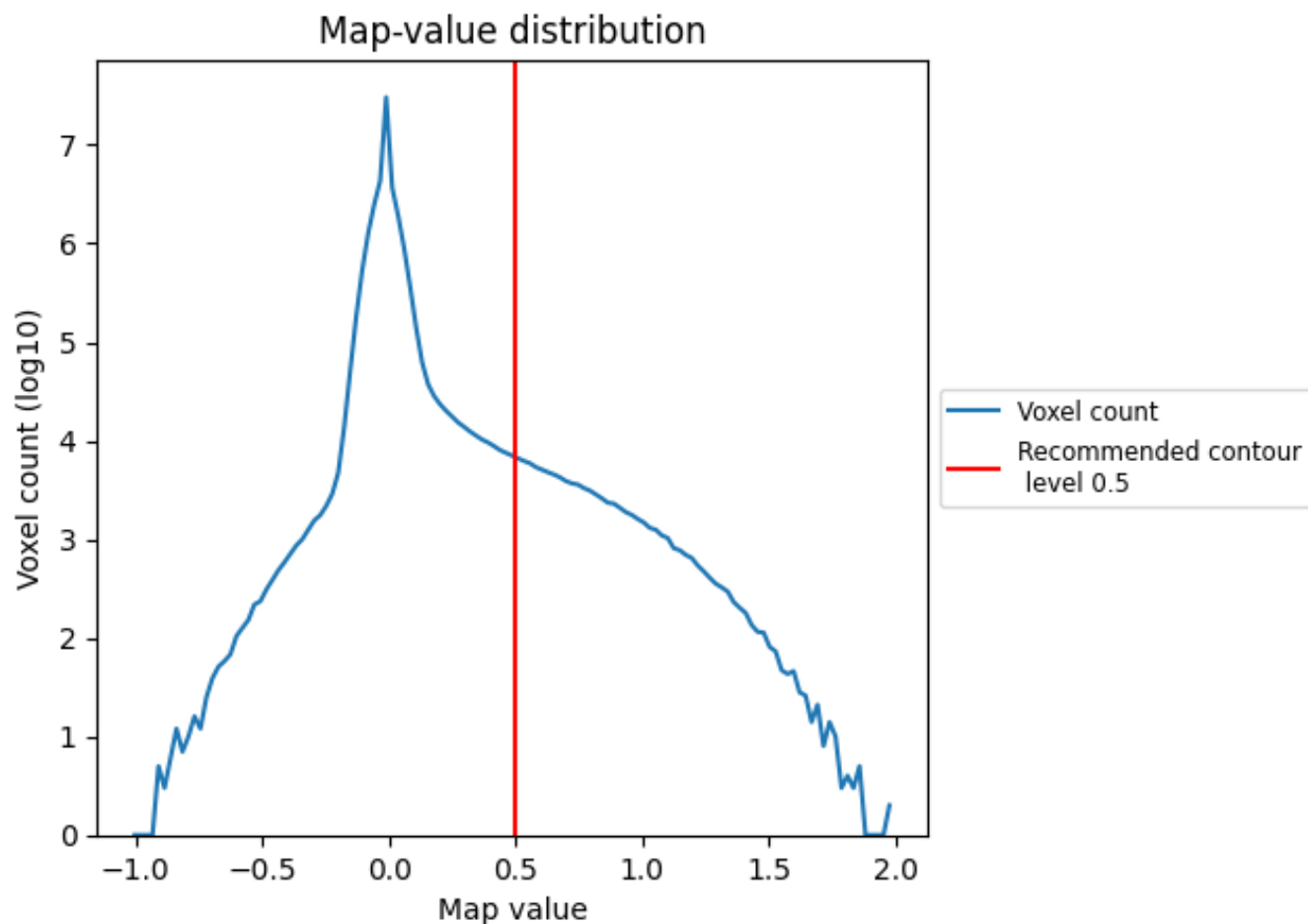
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

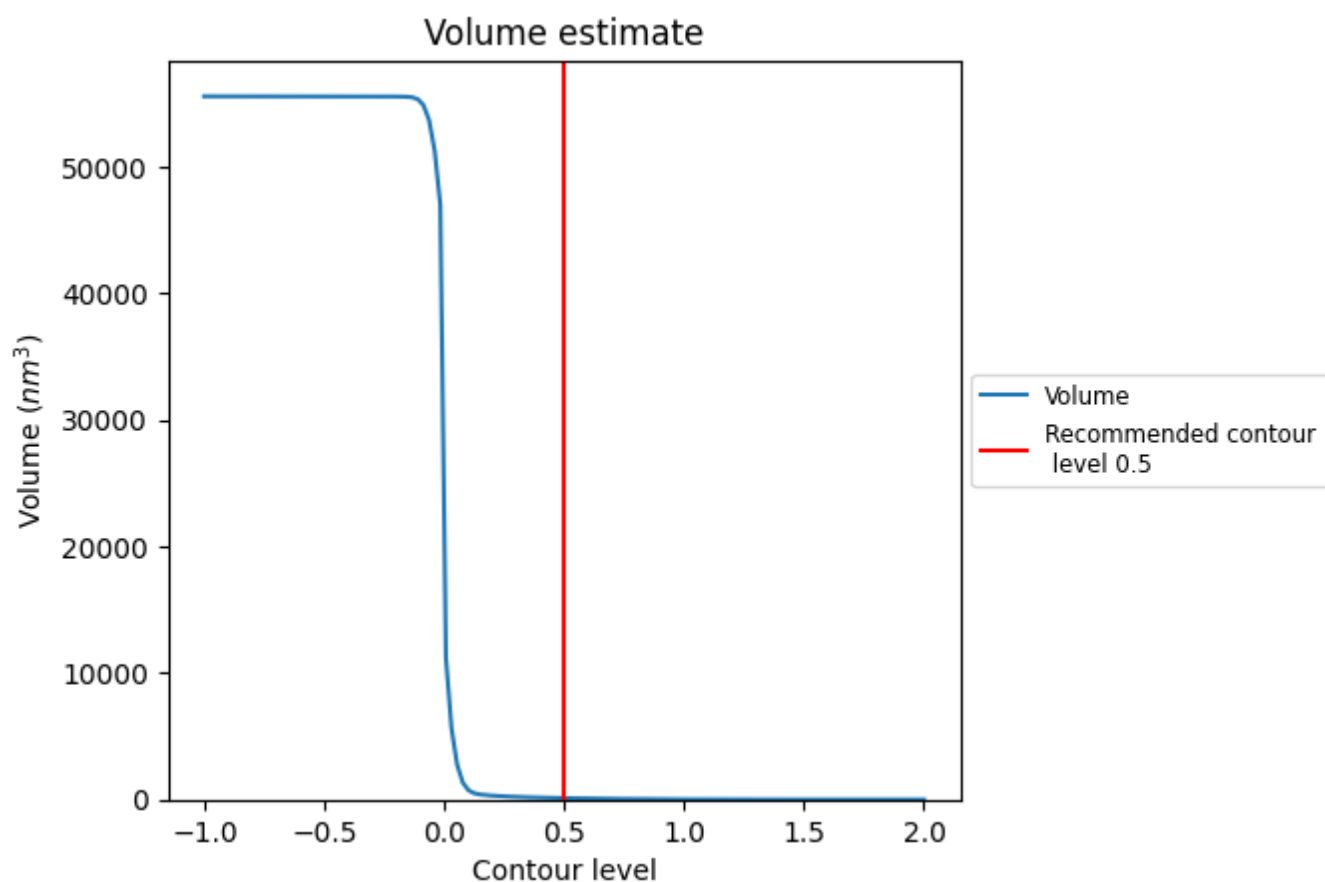
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

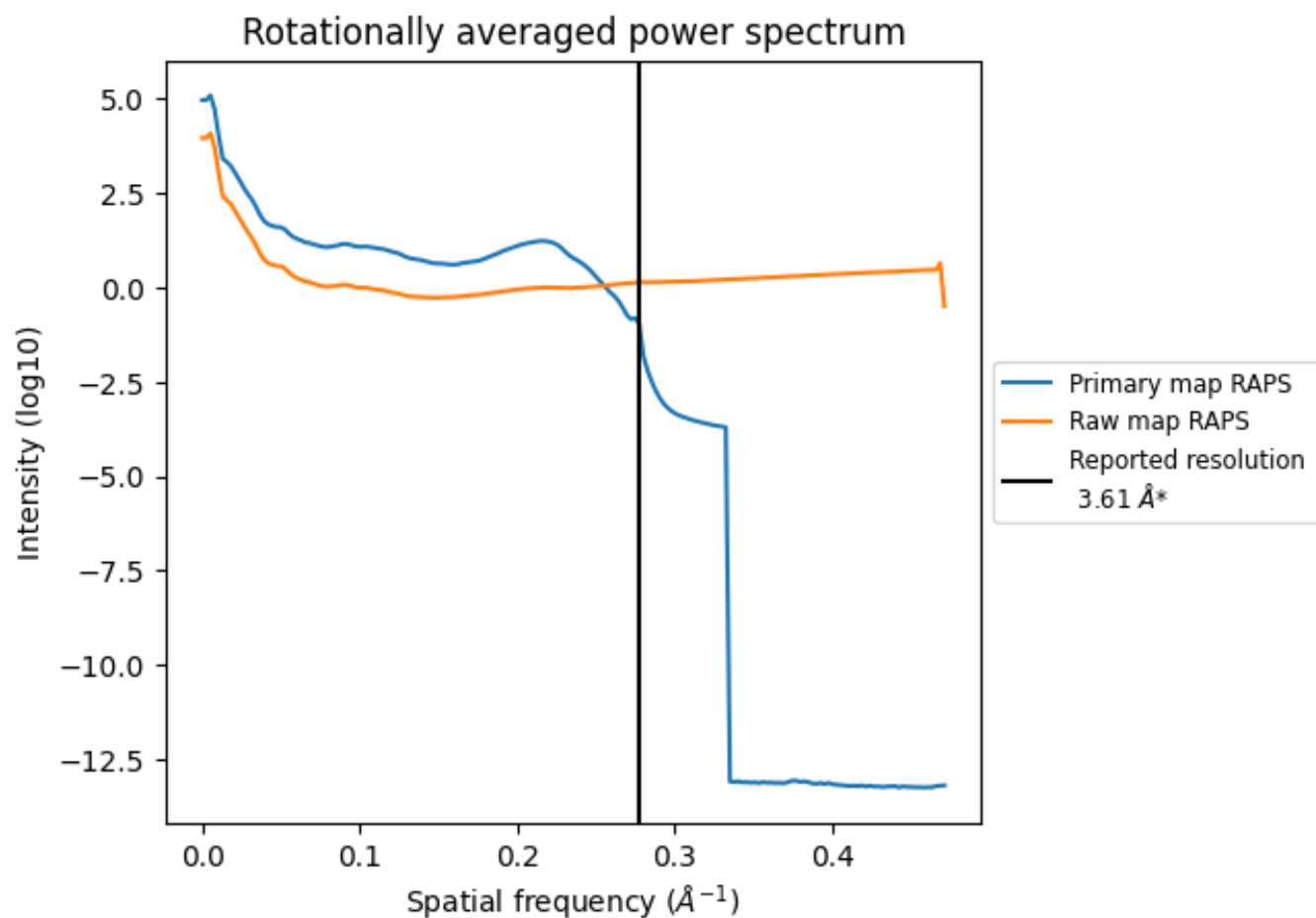
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 111 nm³; this corresponds to an approximate mass of 101 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 ⓘ

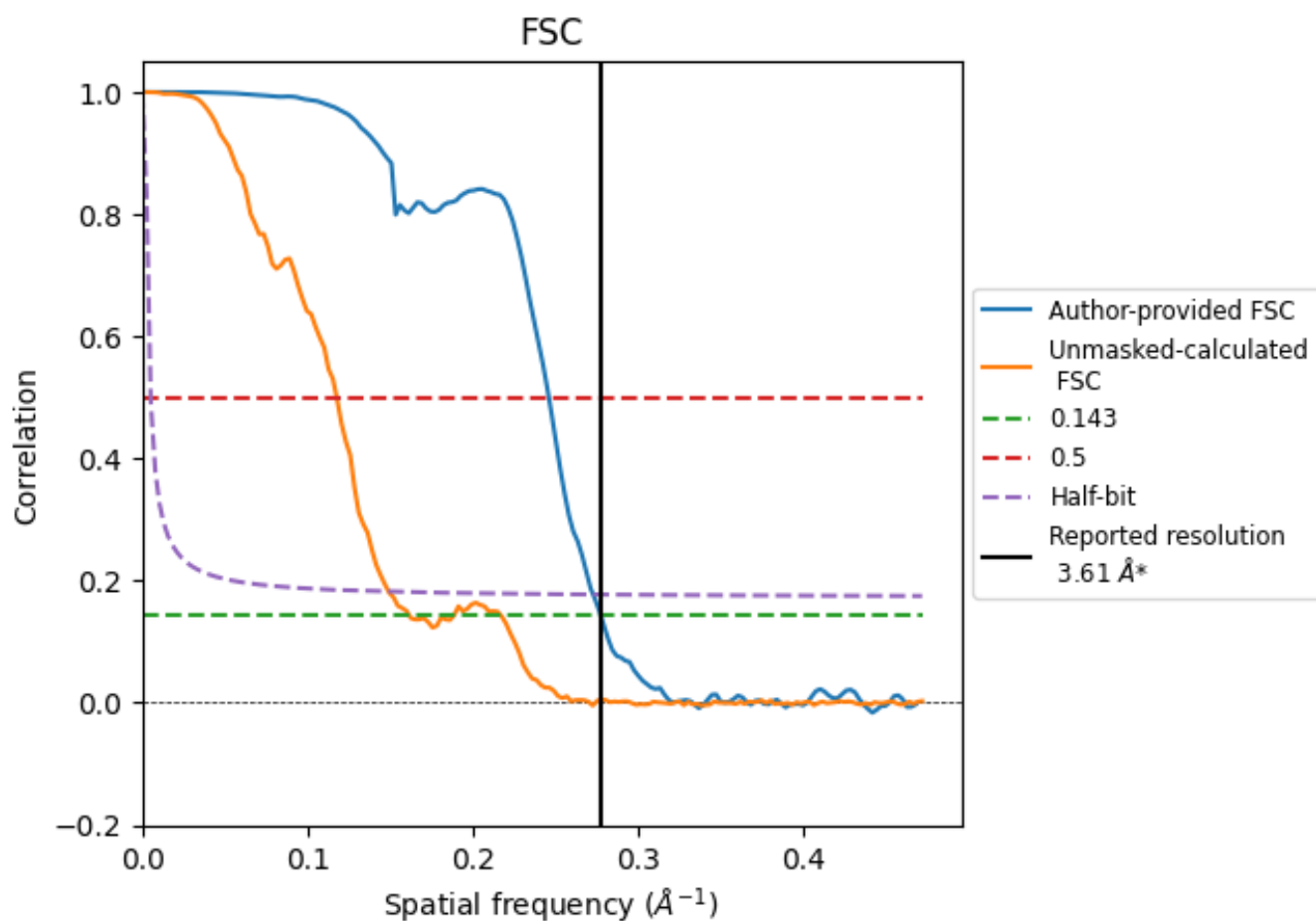


*Reported resolution corresponds to spatial frequency of 0.277 Å⁻¹

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 0.277 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.61	-	-
Author-provided FSC curve	3.61	4.07	3.67
Unmasked-calculated*	6.19	8.50	6.71

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.19 differs from the reported value 3.61 by more than 10 %

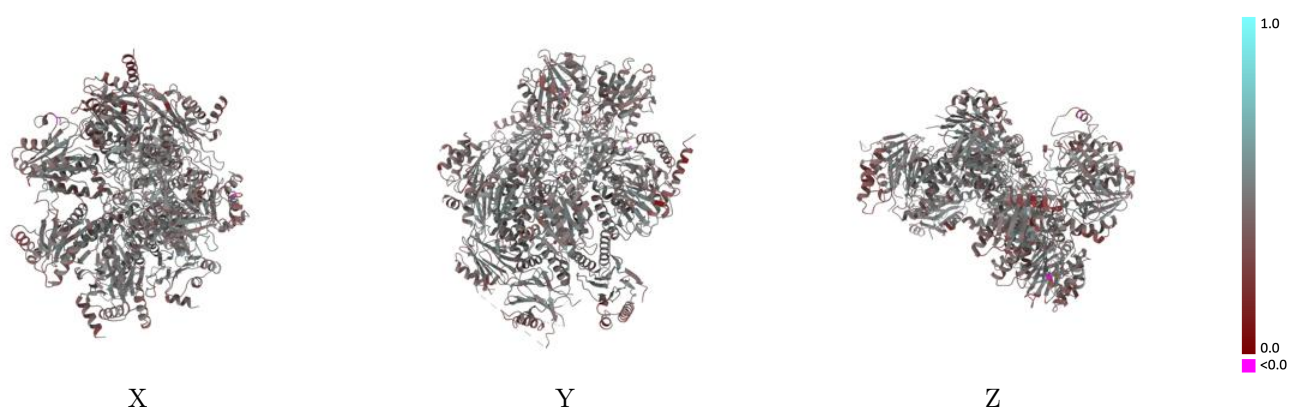
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23508 and PDB model 7LSX. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

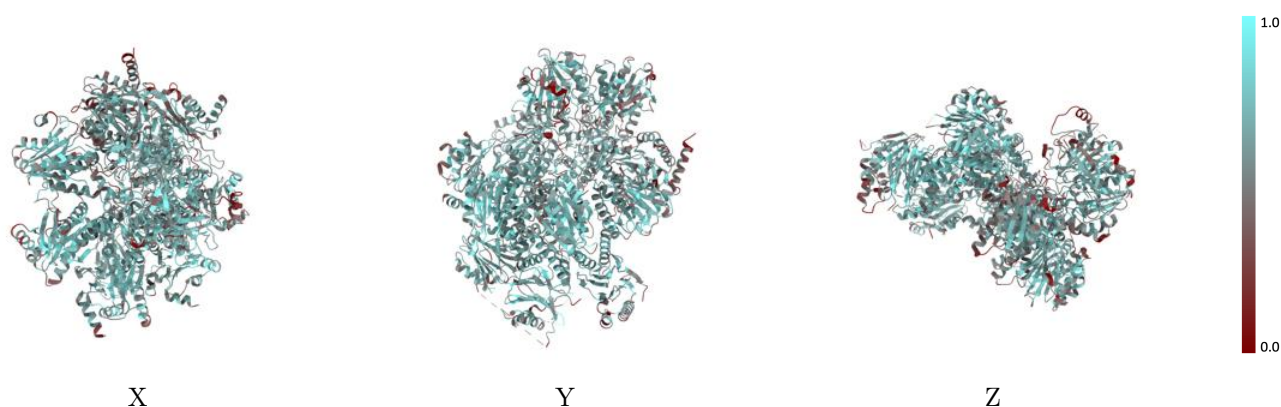
This section was not generated.

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



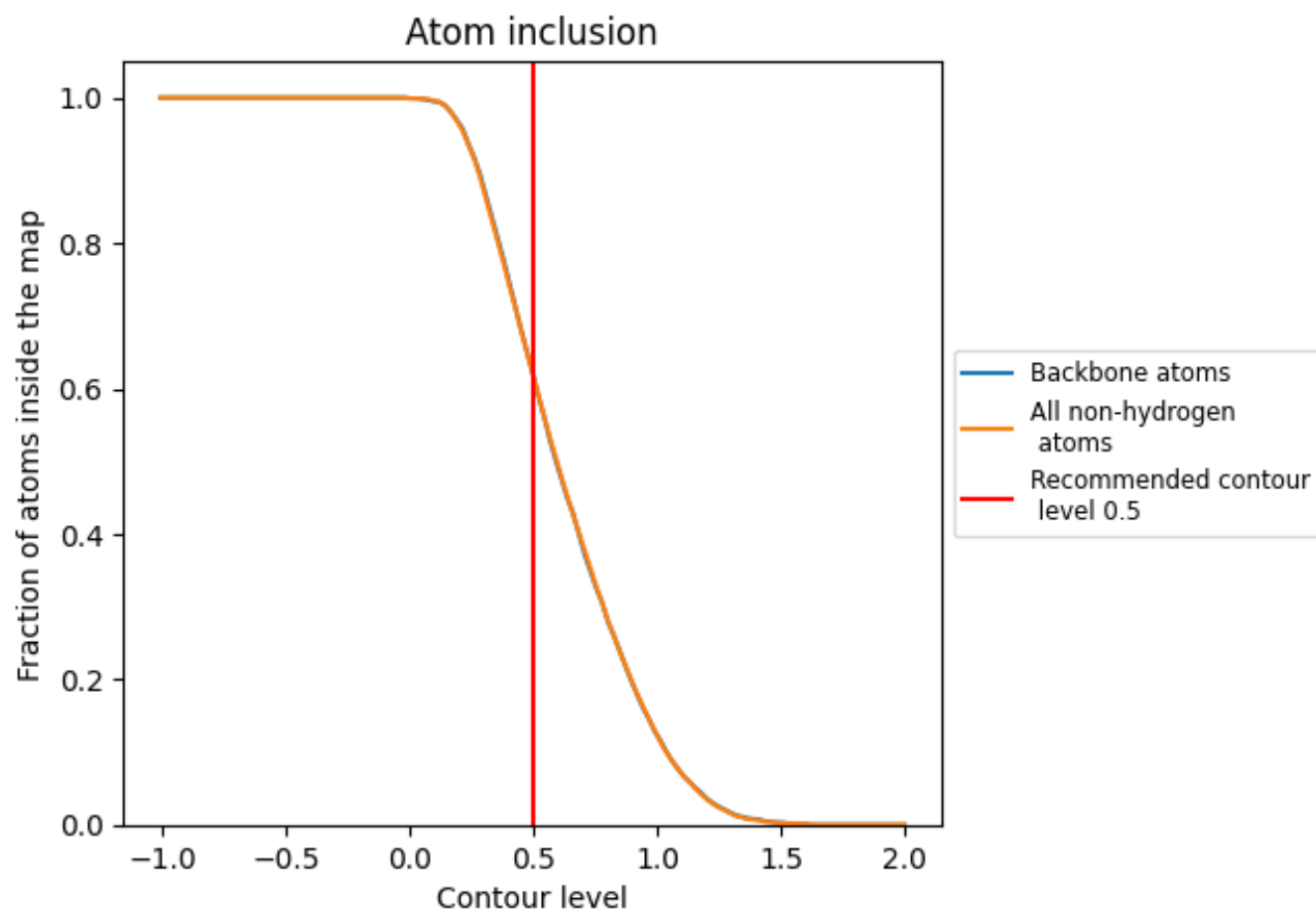
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 62% of all backbone atoms, 62% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6170	0.4410
A	0.6750	0.4500
B	0.6830	0.4620
C	0.6710	0.4560
D	0.6540	0.4490
E	0.5200	0.4160
F	0.6330	0.4360
G	0.6710	0.4470
H	0.5620	0.4480
I	0.6700	0.4380
J	0.6790	0.4490
K	0.6180	0.4230
O	0.5500	0.4320
P	0.5730	0.4330

1.0

0.0

<0.0