



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 07:33 AM UTC

PDB ID : 8TM6 / pdb_00008tm6
EMDB ID : EMD-41380
Title : Human premature 20S proteasome assembly intermediate
Authors : Zhang, H.; Zhao, J.
Deposited on : 2023-07-28
Resolution : 2.80 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:

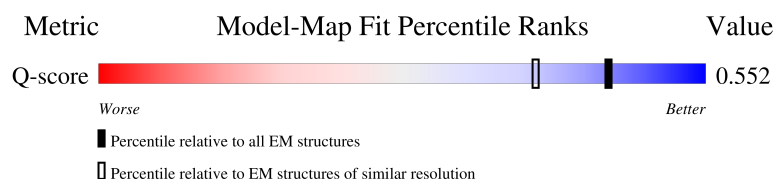
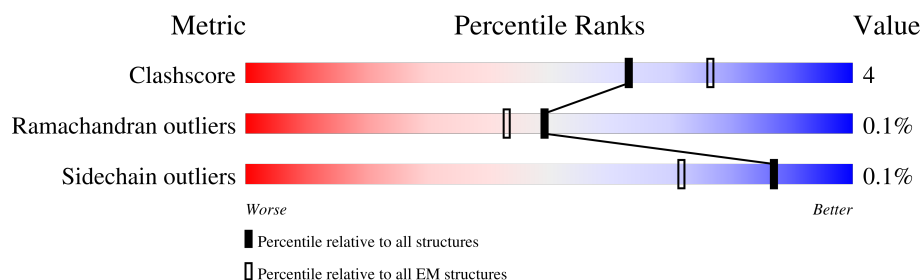
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 2.80 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.



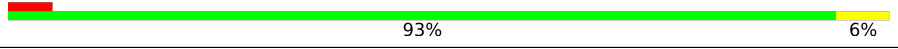
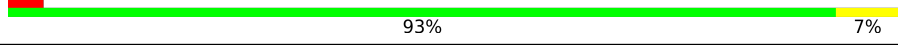
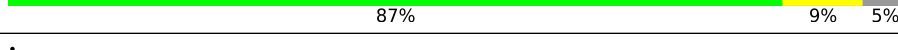
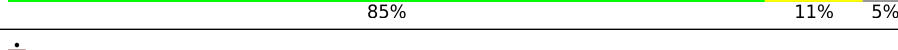
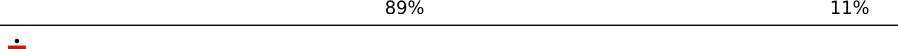
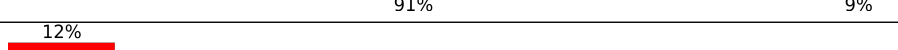
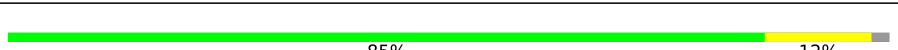
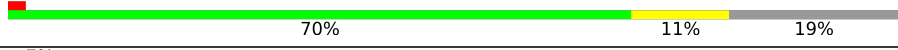
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11806 (2.30 - 3.30)

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	234	 87% 13%
1	O	234	 89% 11%
2	B	261	 5% 89% 7%
2	P	261	 5% 84% 12%

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Mol	Chain	Length	Quality of chain
3	C	248	
3	Q	248	
4	D	241	
4	R	241	
5	E	263	
5	S	263	
6	F	255	
6	T	255	
7	G	246	
7	U	246	
8	H	277	
8	V	277	
9	I	205	
9	W	205	
10	J	201	
10	X	201	
11	K	263	
11	Y	263	
12	L	241	
12	Z	241	
13	M	264	
13	a	264	
14	N	239	
14	b	239	
15	c	288	

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Mol	Chain	Length	Quality of chain
15	f	288	49% 85% 9% 5%
16	d	264	16% 84% 10% 6%
16	g	264	18% 83% 11% 6%
17	e	141	27% 60% 13% 27%
17	h	141	28% 65% 7% 27%

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 58074 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 Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	234	Total	C	N	O	S	0	0
			1780	1143	305	326	6		
1	O	234	Total	C	N	O	S	0	0
			1780	1143	305	326	6		

- Molecule 2 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	251	Total	C	N	O	S	0	0
			1929	1223	337	359	10		
2	P	251	Total	C	N	O	S	0	0
			1929	1223	337	359	10		

- Molecule 3 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	238	Total	C	N	O	S	0	0
			1745	1101	319	320	5		
3	Q	238	Total	C	N	O	S	0	0
			1745	1101	319	320	5		

- Molecule 4 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	241	Total	C	N	O	S	0	0
			1800	1140	303	345	12		
4	R	241	Total	C	N	O	S	0	0
			1800	1140	303	345	12		

- Molecule 5 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	239	Total	C	N	O	S	0	0
			1830	1155	336	328	11		
5	S	239	Total	C	N	O	S	0	0
			1830	1155	336	328	11		

- Molecule 6 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	243	Total	C	N	O	S	0	0
			1847	1178	318	340	11		
6	T	243	Total	C	N	O	S	0	0
			1847	1178	318	340	11		

- Molecule 7 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	245	Total	C	N	O	S	0	0
			1839	1172	315	340	12		
7	U	245	Total	C	N	O	S	0	0
			1839	1172	315	340	12		

- Molecule 8 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	259	Total	C	N	O	S	0	0
			1910	1207	329	362	12		
8	V	259	Total	C	N	O	S	0	0
			1910	1207	329	362	12		

- Molecule 9 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	205	Total	C	N	O	S	0	0
			1575	1005	264	287	19		
9	W	205	Total	C	N	O	S	0	0
			1575	1005	264	287	19		

- Molecule 10 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	197	Total	C	N	O	S	0	0
			1564	1006	265	284	9		

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Mol	Chain	Residues	Atoms					AltConf	Trace
10	X	197	Total	C	N	O	S	0	0
			1564	1006	265	284	9		

- Molecule 11 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	200	Total	C	N	O	S	0	0
			1539	972	271	287	9		
11	Y	200	Total	C	N	O	S	0	0
			1539	972	271	287	9		

- Molecule 12 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	213	Total	C	N	O	S	0	0
			1630	1037	281	302	10		
12	Z	213	Total	C	N	O	S	0	0
			1630	1037	281	302	10		

- Molecule 13 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	215	Total	C	N	O	S	0	0
			1642	1045	288	297	12		
13	a	215	Total	C	N	O	S	0	0
			1642	1045	288	297	12		

- Molecule 14 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	215	Total	C	N	O	S	0	0
			1606	1006	275	313	12		
14	b	215	Total	C	N	O	S	0	0
			1606	1006	275	313	12		

- Molecule 15 is a protein called Proteasome assembly chaperone 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	c	273	Total	C	N	O	S	0	0
			2081	1336	347	381	17		
15	f	273	Total	C	N	O	S	0	0
			2081	1336	347	381	17		

- Molecule 16 is a protein called Proteasome assembly chaperone 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	d	249	Total	C	N	O	S	0	0
			1907	1231	312	352	12		
16	g	249	Total	C	N	O	S	0	0
			1907	1231	312	352	12		

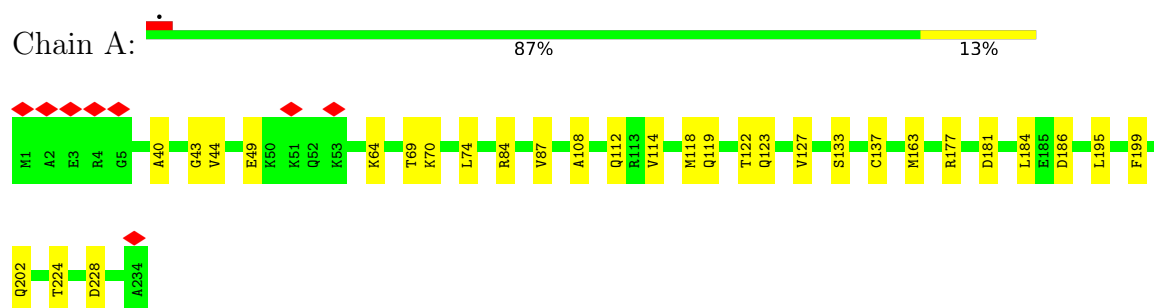
- Molecule 17 is a protein called Proteasome maturation protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	e	103	Total	C	N	O	S	0	0
			813	516	137	156	4		
17	h	103	Total	C	N	O	S	0	0
			813	516	137	156	4		

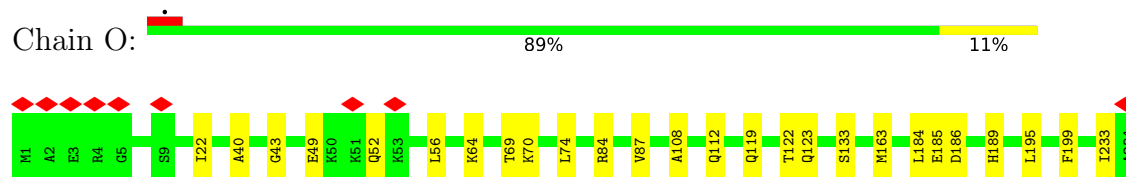
3 Residue-property plots [i](#)

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.

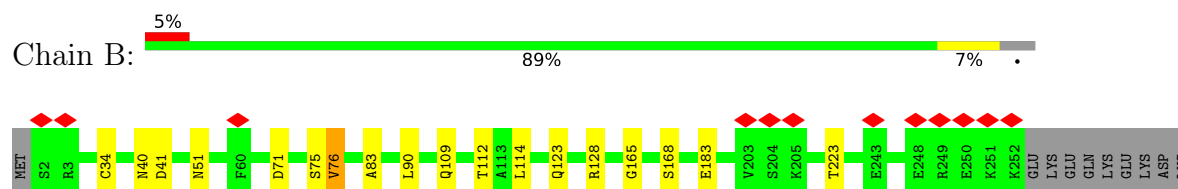
- Molecule 1: Proteasome subunit alpha type-2



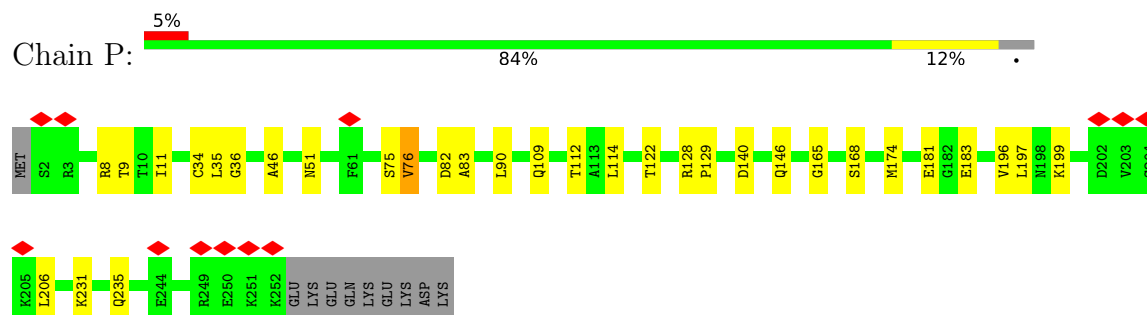
- Molecule 1: Proteasome subunit alpha type-2



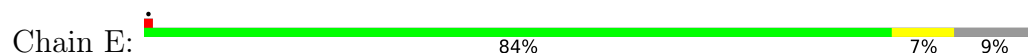
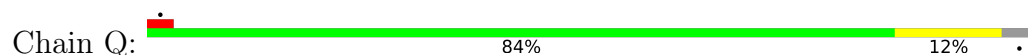
- Molecule 2: Proteasome subunit alpha type-4



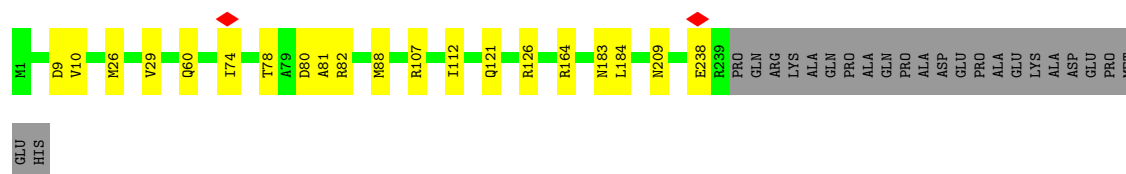
- Molecule 2: Proteasome subunit alpha type-4



Chain C: 84% 12%



Chain S:  83% 8% 9%



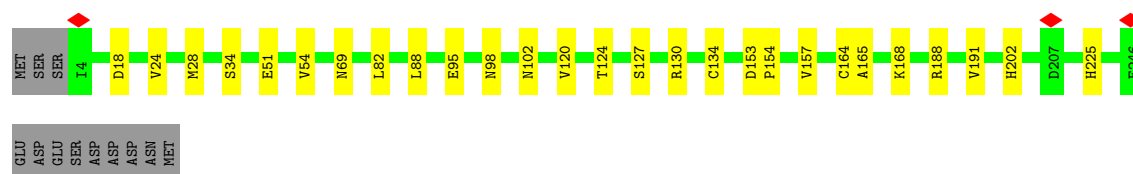
- Molecule 6: Proteasome subunit alpha type-3

Chain F:  87% 9% 5%



- Molecule 6: Proteasome subunit alpha type-3

Chain T:  85% 11% 5%



- Molecule 7: Proteasome subunit alpha type-6

Chain G:  89% 11%



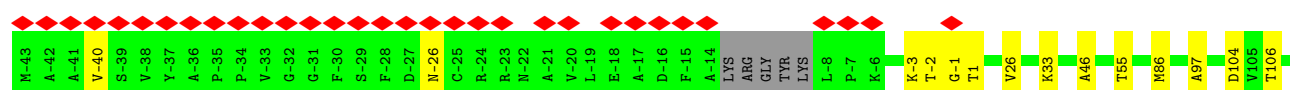
- Molecule 7: Proteasome subunit alpha type-6

Chain U:  91% 9%



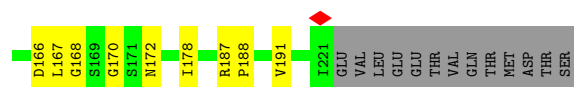
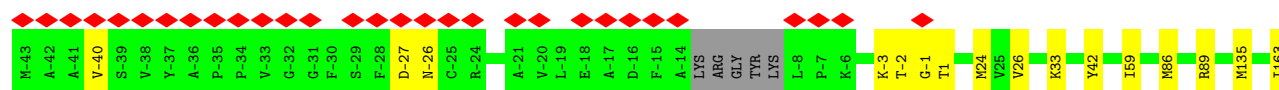
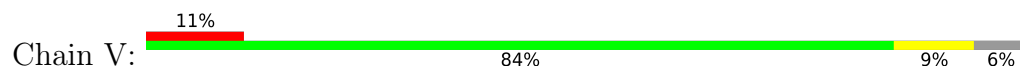
- Molecule 8: Proteasome subunit beta type-7

Chain H:  12% 84% 9% 6%





- Molecule 8: Proteasome subunit beta type-7



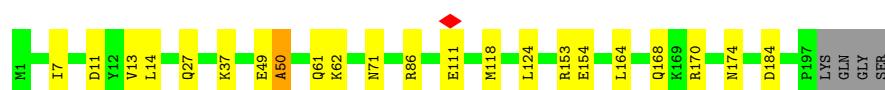
- Molecule 9: Proteasome subunit beta type-3



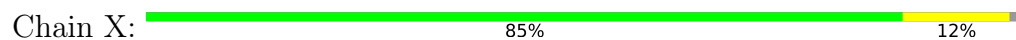
- Molecule 9: Proteasome subunit beta type-3



- Molecule 10: Proteasome subunit beta type-2



- Molecule 10: Proteasome subunit beta type-2



- Molecule 11: Proteasome subunit beta type-5





• Molecule 11: Proteasome subunit beta type-5

Chain Y: 74% 24%



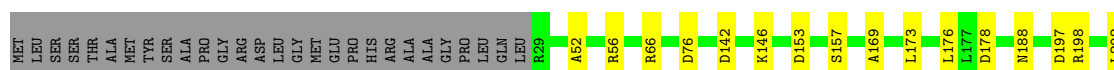
• Molecule 12: Proteasome subunit beta type-1

Chain L: 80% 8% 12%



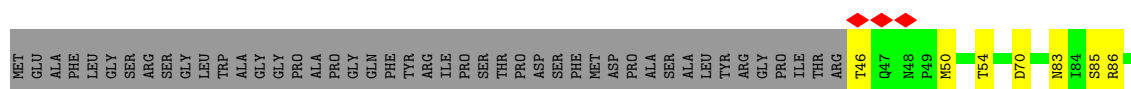
• Molecule 12: Proteasome subunit beta type-1

Chain Z: 79% 9% 12%



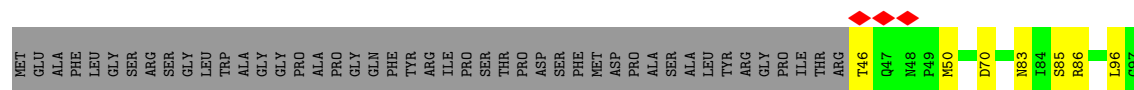
• Molecule 13: Proteasome subunit beta type-4

Chain M: 71% 11% 19%

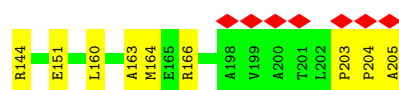
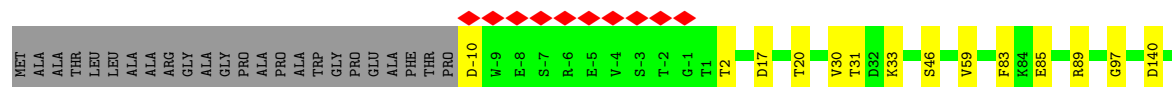
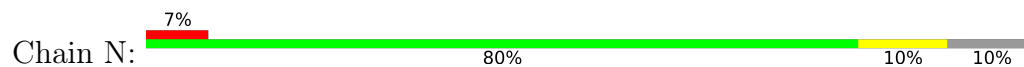


• Molecule 13: Proteasome subunit beta type-4

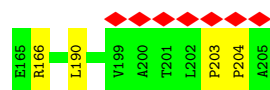
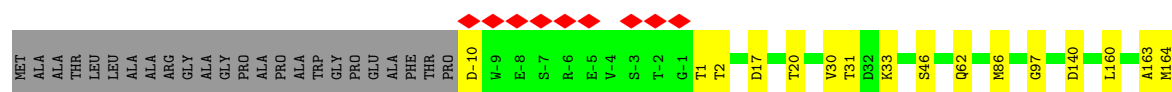
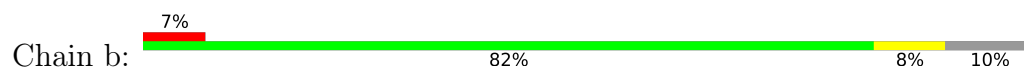
Chain a: 70% 11% 19%



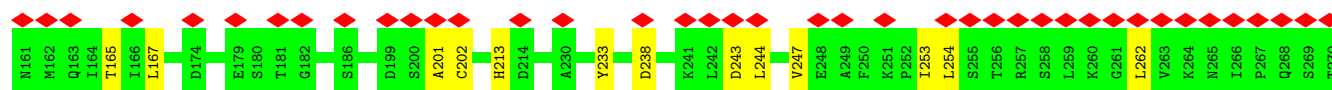
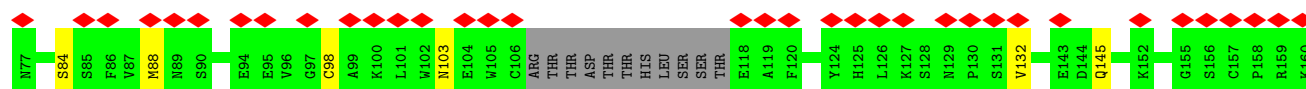
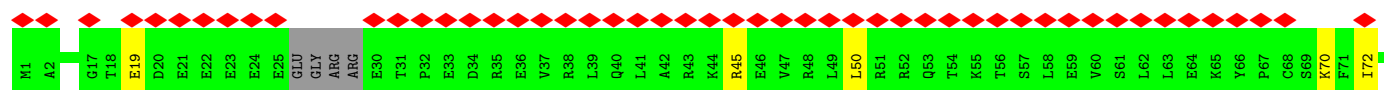
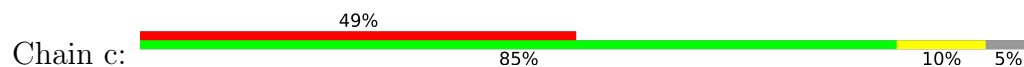
• Molecule 14: Proteasome subunit beta type-6



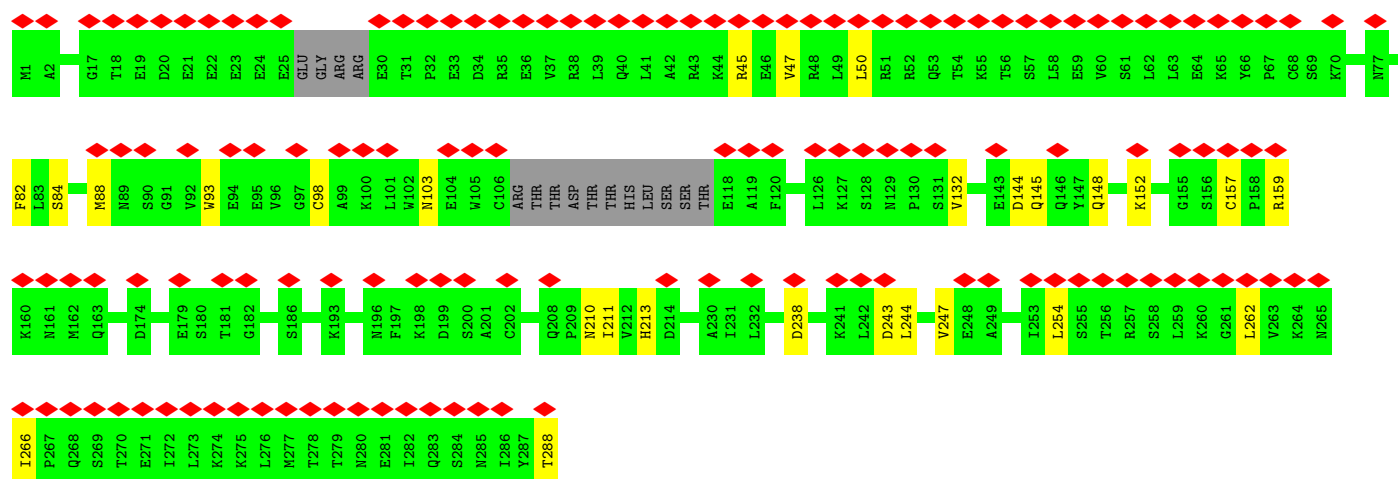
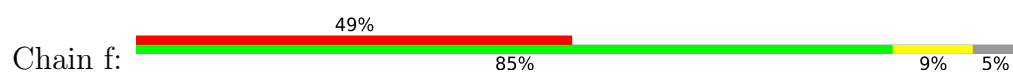
• Molecule 14: Proteasome subunit beta type-6



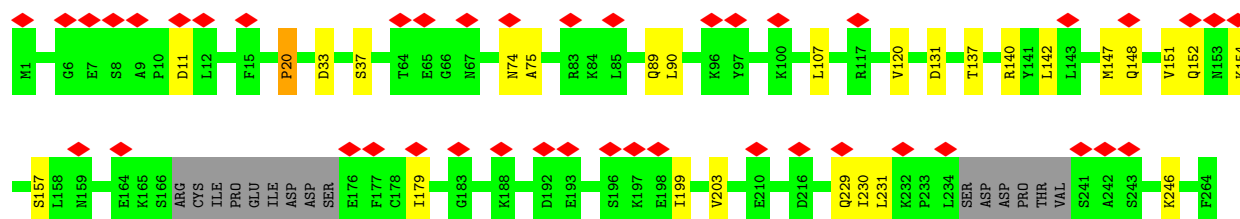
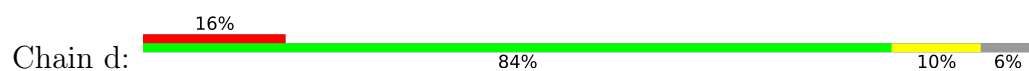
• Molecule 15: Proteasome assembly chaperone 1



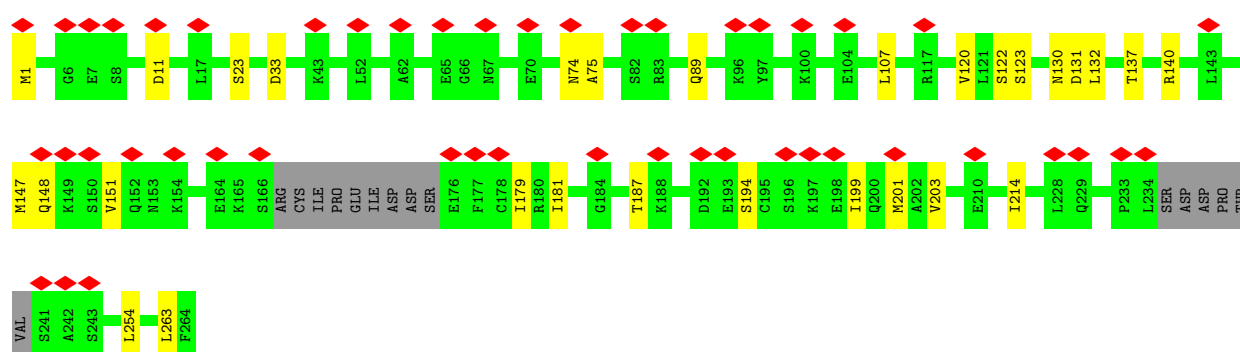
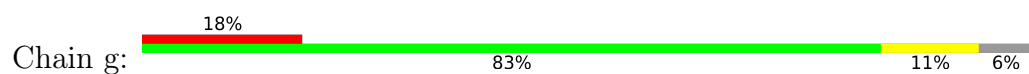
• Molecule 15: Proteasome assembly chaperone 1



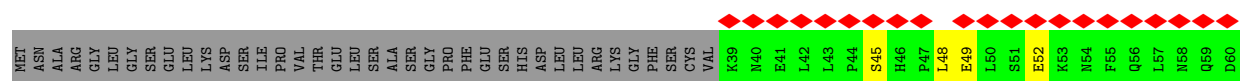
• Molecule 16: Proteasome assembly chaperone 2

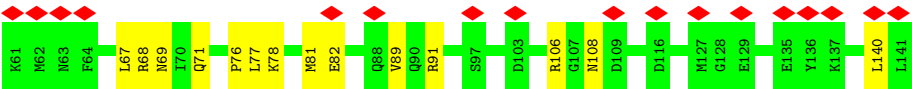


• Molecule 16: Proteasome assembly chaperone 2

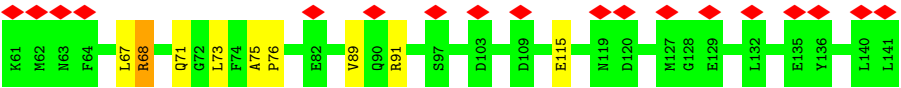
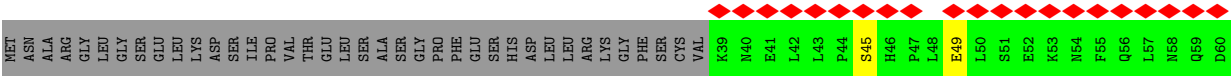


• Molecule 17: Proteasome maturation protein





• Molecule 17: Proteasome maturation protein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	115008	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 ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	1700	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	30.717	Depositor
Minimum map value	-14.922	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	6	Depositor
Map size (Å)	423.99997, 423.99997, 423.99997	wwPDB
Map dimensions	400, 400, 400	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	0.11	0/1819	0.24	0/2470
1	O	0.12	0/1819	0.27	0/2470
2	B	0.10	0/1959	0.23	0/2646
2	P	0.10	0/1959	0.24	0/2646
3	C	0.10	0/1770	0.24	0/2407
3	Q	0.10	0/1770	0.25	0/2407
4	D	0.10	0/1828	0.25	0/2472
4	R	0.11	0/1828	0.26	0/2472
5	E	0.10	0/1865	0.25	0/2525
5	S	0.10	0/1865	0.24	0/2525
6	F	0.12	0/1882	0.25	0/2543
6	T	0.11	0/1882	0.28	0/2543
7	G	0.12	0/1871	0.27	0/2535
7	U	0.13	0/1871	0.30	0/2535
8	H	0.11	0/1942	0.28	0/2635
8	V	0.12	0/1942	0.29	0/2635
9	I	0.11	0/1604	0.27	0/2166
9	W	0.12	0/1604	0.28	0/2166
10	J	0.11	0/1597	0.26	0/2163
10	X	0.11	0/1597	0.28	0/2163
11	K	0.10	0/1570	0.23	0/2125
11	Y	0.10	0/1570	0.23	0/2125
12	L	0.10	0/1660	0.24	0/2240
12	Z	0.10	0/1660	0.24	0/2240
13	M	0.11	0/1675	0.26	0/2270
13	a	0.11	0/1675	0.26	0/2270
14	N	0.12	0/1636	0.25	0/2218
14	b	0.12	0/1636	0.26	0/2218
15	c	0.11	0/2125	0.27	0/2893
15	f	0.10	0/2125	0.26	0/2893
16	d	0.17	0/1948	0.39	2/2645 (0.1%)
16	g	0.12	0/1948	0.30	0/2645
17	e	0.13	0/827	0.35	0/1117
17	h	0.09	0/827	0.27	0/1117

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.11	0/59156	0.27	2/80140 (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
7	U	0	1
17	h	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	d	20	PRO	CA-N-CD	-9.90	98.14	112.00
16	d	20	PRO	N-CD-CG	-5.33	95.21	103.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
7	U	153	LYS	Peptide
17	h	68	ARG	Sidechain

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	1780	0	1746	20	0
1	O	1780	0	1746	19	0
2	B	1929	0	1902	12	0
2	P	1929	0	1902	22	0
3	C	1745	0	1663	22	0
3	Q	1745	0	1663	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1800	0	1757	12	0
4	R	1800	0	1757	14	0
5	E	1830	0	1793	16	0
5	S	1830	0	1793	18	0
6	F	1847	0	1794	16	0
6	T	1847	0	1794	19	0
7	G	1839	0	1792	19	0
7	U	1839	0	1792	14	0
8	H	1910	0	1913	17	0
8	V	1910	0	1913	20	0
9	I	1575	0	1582	6	0
9	W	1575	0	1582	14	0
10	J	1564	0	1559	18	0
10	X	1564	0	1559	21	0
11	K	1539	0	1482	8	0
11	Y	1539	0	1482	4	0
12	L	1630	0	1622	14	0
12	Z	1630	0	1622	16	0
13	M	1642	0	1613	18	0
13	a	1642	0	1613	19	0
14	N	1606	0	1557	14	0
14	b	1606	0	1557	13	0
15	c	2081	0	2014	20	0
15	f	2081	0	2014	19	0
16	d	1907	0	1907	17	0
16	g	1907	0	1907	22	0
17	e	813	0	801	16	0
17	h	813	0	801	12	0
All	All	58074	0	56994	486	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 4.

All (486) 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)
13:M:224:ARG:NH1	8:V:135:MET:SD	2.40	0.94
8:H:135:MET:SD	13:a:224:ARG:NH1	2.46	0.89
4:D:52:LYS:NZ	4:D:64:ILE:O	2.11	0.82
16:g:147:MET:HE3	16:g:151:VAL:HG21	1.63	0.81
4:R:52:LYS:NZ	4:R:64:ILE:O	2.14	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:V:-26:ASN:ND2	9:W:101:GLY:O	2.16	0.79
8:H:-26:ASN:ND2	9:I:101:GLY:O	2.17	0.78
5:S:183:ASN:OD1	5:S:184:LEU:N	2.19	0.75
16:g:33:ASP:OD1	16:g:89:GLN:NE2	2.21	0.74
9:W:178:ASP:OD2	9:W:181:SER:OG	2.04	0.73
16:d:33:ASP:OD1	16:d:89:GLN:NE2	2.20	0.73
2:P:140:ASP:OD1	2:P:146:GLN:NE2	2.23	0.72
5:S:164:ARG:NH2	15:f:243:ASP:OD2	2.22	0.72
9:I:177:ARG:NH1	12:Z:178:ASP:OD2	2.23	0.71
15:c:45:ARG:O	15:c:145:GLN:NE2	2.23	0.71
16:d:74:ASN:OD1	16:d:75:ALA:N	2.24	0.70
9:I:65:GLN:OE1	10:J:86:ARG:NH2	2.24	0.70
15:f:45:ARG:O	15:f:145:GLN:NE2	2.23	0.70
8:H:-3:LYS:O	8:H:-2:THR:OG1	2.09	0.70
6:T:120:VAL:O	6:T:124:THR:HG23	1.92	0.69
17:e:106:ARG:NH1	17:e:108:ASN:OD1	2.25	0.69
6:F:98:ASN:HB2	17:e:68:ARG:HH12	1.57	0.69
4:D:160:GLY:O	5:E:82:ARG:NH2	2.25	0.69
6:F:120:VAL:O	6:F:124:THR:HG23	1.92	0.68
4:R:160:GLY:O	5:S:82:ARG:NH2	2.25	0.68
8:V:-3:LYS:O	8:V:-2:THR:OG1	2.11	0.68
12:L:213:ARG:NH1	8:V:26:VAL:O	2.28	0.67
14:N:85:GLU:OE2	14:N:89:ARG:NH2	2.27	0.67
14:b:160:LEU:O	14:b:164:MET:HG3	1.93	0.67
9:W:65:GLN:OE1	10:X:86:ARG:NH2	2.27	0.67
8:V:89:ARG:NH2	17:h:115:GLU:OE2	2.28	0.67
2:P:183:GLU:N	2:P:183:GLU:OE2	2.29	0.66
10:X:154:GLU:OE1	10:X:154:GLU:N	2.27	0.66
17:h:115:GLU:N	17:h:115:GLU:OE1	2.28	0.66
2:B:40:ASN:OD1	2:B:41:ASP:N	2.29	0.66
7:G:192:GLU:OE2	7:G:192:GLU:N	2.29	0.65
12:L:88:ASP:OD1	13:M:142:TYR:OH	2.11	0.64
6:F:164:CYS:SG	6:F:165:ALA:N	2.70	0.64
17:e:67:LEU:O	17:e:71:GLN:N	2.31	0.64
3:Q:178:ASP:OD1	3:Q:179:GLU:N	2.32	0.63
10:X:109:GLU:OE2	10:X:109:GLU:N	2.30	0.63
10:J:153:ARG:NH1	10:J:184:ASP:OD2	2.30	0.63
13:M:50:MET:HE2	13:M:50:MET:HA	1.81	0.63
6:F:98:ASN:OD1	6:F:99:PHE:N	2.31	0.63
14:N:160:LEU:O	14:N:164:MET:HG3	1.99	0.63
2:P:165:GLY:O	2:P:168:SER:OG	2.17	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:f:50:LEU:O	15:f:98:CYS:N	2.31	0.62
5:E:180:MET:HE2	5:E:180:MET:HA	1.80	0.62
10:X:29:LYS:NZ	10:X:31:ASP:OD1	2.31	0.62
16:g:131:ASP:OD1	16:g:132:LEU:N	2.32	0.62
8:H:164:PHE:O	12:Z:66:ARG:NH2	2.33	0.61
5:E:164:ARG:NH1	15:c:238:ASP:OD1	2.34	0.61
15:c:50:LEU:O	15:c:98:CYS:N	2.33	0.61
5:E:60:GLN:OE1	15:c:288:THR:HB	2.00	0.61
5:E:164:ARG:NH2	15:c:243:ASP:OD2	2.33	0.61
10:J:170:ARG:NH1	10:X:27:GLN:O	2.34	0.61
11:K:78:ALA:O	11:K:82:LEU:HD13	2.01	0.61
14:b:46:SER:OG	14:b:97:GLY:O	2.16	0.61
15:f:103:ASN:ND2	15:f:145:GLN:OE1	2.33	0.61
10:J:27:GLN:O	10:X:170:ARG:NH1	2.34	0.60
3:C:42:VAL:HG13	3:C:191:VAL:HG21	1.82	0.60
3:Q:90:GLU:OE2	3:Q:106:TYR:OH	2.14	0.60
1:O:56:LEU:O	1:O:56:LEU:HD12	2.01	0.60
15:f:213:HIS:O	15:f:213:HIS:ND1	2.35	0.60
3:C:159:ASN:OD1	3:C:160:ALA:N	2.36	0.59
3:C:42:VAL:CG1	3:C:191:VAL:HG21	2.33	0.59
15:c:213:HIS:O	15:c:213:HIS:ND1	2.36	0.59
6:T:164:CYS:SG	6:T:165:ALA:N	2.76	0.59
10:J:37:LYS:O	10:J:61:GLN:NE2	2.36	0.58
5:S:9:ASP:OD1	5:S:10:VAL:N	2.35	0.58
6:T:153:ASP:HB3	6:T:157:VAL:HG22	1.84	0.58
13:M:46:THR:N	13:M:150:PRO:O	2.36	0.58
9:W:30:ILE:N	9:W:30:ILE:HD12	2.19	0.58
3:Q:159:ASN:OD1	3:Q:160:ALA:N	2.37	0.58
14:N:46:SER:OG	14:N:97:GLY:O	2.17	0.58
6:T:95:GLU:OE1	17:h:73:LEU:N	2.31	0.58
15:c:103:ASN:ND2	15:c:145:GLN:OE1	2.37	0.58
13:M:70:ASP:OD1	13:M:86:ARG:NH2	2.37	0.57
13:a:46:THR:N	13:a:150:PRO:O	2.38	0.57
5:S:74:ILE:HD12	5:S:81:ALA:HB1	1.86	0.57
5:S:107:ARG:NH2	13:a:119:GLU:OE1	2.30	0.57
16:g:74:ASN:OD1	16:g:75:ALA:N	2.37	0.57
12:L:178:ASP:OD2	9:W:177:ARG:NH1	2.37	0.57
10:X:153:ARG:NH1	10:X:184:ASP:OD2	2.38	0.57
6:F:102:ASN:ND2	17:e:69:ASN:OD1	2.38	0.57
1:O:133:SER:OG	1:O:163:MET:SD	2.62	0.57
6:F:143:VAL:HG13	6:F:144:ASN:OD1	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:97:PHE:O	12:L:94:LYS:NZ	2.34	0.56
9:I:161:ASP:OD1	9:I:162:HIS:N	2.38	0.56
11:K:148:GLU:N	11:K:148:GLU:OE2	2.37	0.56
12:Z:202:LEU:O	12:Z:206:VAL:HG22	2.06	0.56
16:d:147:MET:HE3	16:d:151:VAL:HG21	1.88	0.56
7:G:155:ASP:OD1	7:G:155:ASP:C	2.49	0.55
14:N:-10:ASP:OD1	14:N:-10:ASP:N	2.39	0.55
1:A:133:SER:OG	1:A:163:MET:SD	2.64	0.55
5:S:164:ARG:NH1	15:f:238:ASP:OD1	2.36	0.55
10:X:37:LYS:O	10:X:61:GLN:NE2	2.38	0.55
2:P:35:LEU:HD23	2:P:36:GLY:N	2.22	0.55
10:X:27:GLN:NE2	10:X:174:ASN:OD1	2.40	0.55
17:e:68:ARG:HB2	17:e:76:PRO:HG2	1.89	0.55
16:g:147:MET:CE	16:g:151:VAL:HG21	2.36	0.55
7:U:155:ASP:OD1	7:U:155:ASP:C	2.49	0.54
2:B:51:ASN:OD1	2:B:51:ASN:O	2.25	0.54
12:L:197:ASP:OD1	12:L:198:ARG:N	2.40	0.54
14:N:17:ASP:OD1	14:N:33:LYS:NZ	2.40	0.54
1:O:22:ILE:HD11	1:O:122:THR:OG1	2.07	0.54
16:d:137:THR:O	16:d:140:ARG:NH1	2.39	0.54
5:E:74:ILE:HD12	5:E:81:ALA:HB1	1.90	0.54
13:a:50:MET:HA	13:a:50:MET:HE2	1.90	0.54
16:g:107:LEU:HD22	16:g:199:ILE:HD11	1.90	0.54
16:g:122:SER:OG	16:g:123:SER:N	2.40	0.54
16:g:179:ILE:HG23	16:g:179:ILE:O	2.06	0.54
14:b:-10:ASP:OD1	14:b:-10:ASP:N	2.40	0.53
15:f:157:CYS:O	15:f:159:ARG:NH1	2.40	0.53
7:G:115:CYS:SG	7:G:160:TYR:HB2	2.49	0.53
10:J:174:ASN:OD1	10:J:174:ASN:O	2.26	0.53
17:e:77:LEU:HD21	17:e:81:MET:HE3	1.90	0.53
1:A:123:GLN:NE2	2:B:128:ARG:O	2.42	0.53
8:H:187:ARG:HB3	8:H:188:PRO:HD3	1.91	0.53
14:b:190:LEU:C	14:b:190:LEU:HD12	2.33	0.53
2:P:51:ASN:OD1	2:P:51:ASN:O	2.27	0.53
10:X:158:GLU:HA	10:X:158:GLU:OE2	2.08	0.53
13:M:83:ASN:OD1	13:M:231:ARG:NH2	2.42	0.52
7:U:72:ILE:HG21	7:U:114:LEU:HD21	1.91	0.52
1:O:184:LEU:C	1:O:185:GLU:OE1	2.52	0.52
14:b:17:ASP:OD1	14:b:33:LYS:NZ	2.42	0.52
15:c:254:LEU:C	15:c:254:LEU:HD12	2.35	0.52
1:O:56:LEU:HD13	7:U:165:ALA:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:234:LEU:O	4:R:238:ILE:HG13	2.10	0.52
16:g:194:SER:CB	16:g:201:MET:HE2	2.39	0.52
3:Q:7:ILE:HD12	3:Q:7:ILE:H	1.74	0.52
12:Z:197:ASP:OD1	12:Z:198:ARG:N	2.43	0.52
13:a:85:SER:OG	13:a:233:GLN:OE1	2.25	0.52
3:Q:120:GLN:OE1	4:R:133:MET:HE1	2.08	0.52
8:V:24:MET:HE3	8:V:24:MET:HA	1.92	0.52
16:d:131:ASP:OD1	16:d:131:ASP:N	2.42	0.52
16:g:137:THR:O	16:g:140:ARG:NH1	2.43	0.52
8:V:187:ARG:HB3	8:V:188:PRO:HD3	1.91	0.52
15:f:47:VAL:O	15:f:152:LYS:NZ	2.38	0.52
7:G:72:ILE:HG21	7:G:114:LEU:HD21	1.91	0.52
13:M:85:SER:OG	13:M:233:GLN:OE1	2.27	0.52
9:W:71:LEU:HD11	9:W:82:ILE:HG21	1.92	0.52
13:a:70:ASP:OD1	13:a:86:ARG:NH2	2.39	0.52
8:H:46:ALA:HB3	8:H:97:ALA:HB3	1.92	0.52
1:O:123:GLN:NE2	2:P:128:ARG:O	2.42	0.52
11:K:82:LEU:O	11:K:86:MET:HG3	2.09	0.51
13:M:211:ARG:NH1	13:M:245:GLU:OE2	2.36	0.51
3:C:145:TYR:CE2	3:C:155:ALA:HB2	2.45	0.51
9:I:71:LEU:HD11	9:I:82:ILE:HG21	1.93	0.51
9:W:135:ASP:OD1	9:W:136:PHE:N	2.36	0.51
16:g:23:SER:OG	16:g:123:SER:N	2.43	0.51
12:L:202:LEU:O	12:L:206:VAL:HG22	2.11	0.51
3:Q:178:ASP:OD1	3:Q:178:ASP:C	2.53	0.51
15:c:167:LEU:HD22	15:c:253:ILE:HD13	1.92	0.51
1:A:84:ARG:NH1	7:G:120:ASP:OD1	2.43	0.51
1:A:202:GLN:OE1	1:A:202:GLN:N	2.35	0.51
2:P:76:VAL:HG21	2:P:83:ALA:HB1	1.93	0.51
13:a:114:GLN:NE2	13:a:118:ASP:OD1	2.43	0.51
13:M:253:ASN:ND2	13:M:255:ASP:OD2	2.44	0.51
6:F:88:LEU:HD12	6:F:134:CYS:SG	2.51	0.51
6:F:98:ASN:OD1	6:F:98:ASN:C	2.53	0.50
2:B:76:VAL:HG21	2:B:83:ALA:HB1	1.94	0.50
14:N:30:VAL:HG11	13:a:256:ILE:HG13	1.94	0.50
10:X:174:ASN:OD1	10:X:174:ASN:O	2.30	0.50
16:g:1:MET:SD	16:g:1:MET:N	2.73	0.50
13:M:86:ARG:NH1	13:M:98:ALA:O	2.44	0.50
1:A:224:THR:O	1:A:228:ASP:OD1	2.30	0.50
3:Q:209:ALA:HB1	3:Q:217:LEU:HD11	1.93	0.50
3:C:155:ALA:HB3	4:D:63:SER:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:166:ARG:NH1	9:W:34:MET:O	2.45	0.50
1:O:84:ARG:NH1	7:U:120:ASP:OD1	2.45	0.50
15:c:132:VAL:HG11	15:c:262:LEU:HD11	1.94	0.50
1:A:108:ALA:O	1:A:112:GLN:HG3	2.12	0.50
12:L:46:GLU:OE2	12:L:46:GLU:HA	2.12	0.49
5:S:26:MET:HE1	16:g:263:LEU:HD11	1.94	0.49
16:g:130:ASN:OD1	16:g:131:ASP:N	2.41	0.49
17:h:89:VAL:O	17:h:89:VAL:HG13	2.12	0.49
10:J:111:GLU:HA	10:J:111:GLU:OE2	2.12	0.49
13:M:256:ILE:HG13	14:b:30:VAL:HG11	1.93	0.49
6:T:88:LEU:HD12	6:T:134:CYS:SG	2.53	0.49
6:F:98:ASN:HB2	17:e:68:ARG:NH1	2.26	0.49
15:f:254:LEU:C	15:f:254:LEU:HD12	2.37	0.49
11:K:14:VAL:HG21	11:K:42:LEU:HD11	1.94	0.49
10:X:18:ASP:C	10:X:18:ASP:OD1	2.56	0.49
2:B:109:GLN:NE2	10:J:71:ASN:OD1	2.46	0.49
10:X:111:GLU:OE1	10:X:111:GLU:HA	2.13	0.49
5:S:74:ILE:CD1	5:S:81:ALA:HB1	2.42	0.49
9:W:73:LEU:O	9:W:77:LYS:HG3	2.13	0.49
15:c:275:LYS:O	15:c:279:THR:HG23	2.13	0.49
3:C:209:ALA:HB1	3:C:217:LEU:HD21	1.94	0.48
7:U:39:SER:HG	7:U:81:THR:HG1	1.60	0.48
3:C:140:GLY:O	3:C:213:ARG:NH2	2.46	0.48
5:E:107:ARG:NE	13:M:124:ASP:OD2	2.36	0.48
2:P:181:GLU:OE1	2:P:181:GLU:N	2.31	0.48
3:Q:116:GLN:O	3:Q:119:THR:OG1	2.28	0.48
4:R:202:LEU:O	4:R:206:MET:HG3	2.13	0.48
7:U:203:SER:OG	7:U:208:ILE:O	2.29	0.48
7:G:187:PHE:CD1	7:G:187:PHE:N	2.82	0.48
1:O:189:HIS:HB3	1:O:233:ILE:HD13	1.95	0.48
13:a:164:GLU:OE2	13:a:164:GLU:HA	2.13	0.48
2:P:34:CYS:SG	2:P:75:SER:OG	2.50	0.48
4:D:202:LEU:O	4:D:206:MET:HG3	2.12	0.48
8:H:97:ALA:HB1	8:H:127:MET:SD	2.53	0.48
1:O:119:GLN:OE1	17:h:91:ARG:NH2	2.46	0.48
16:g:11:ASP:C	16:g:11:ASP:OD1	2.56	0.48
2:B:71:ASP:OD1	2:B:223:THR:OG1	2.28	0.48
2:P:109:GLN:NE2	10:X:71:ASN:OD1	2.47	0.48
11:K:167:ASP:OD1	11:K:169:TYR:N	2.44	0.48
1:O:184:LEU:C	1:O:186:ASP:H	2.21	0.48
4:R:236:GLU:OE1	4:R:236:GLU:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:188:ILE:O	3:C:192:ILE:HG12	2.13	0.48
6:F:127:SER:O	7:G:3:ARG:NH2	2.46	0.48
5:S:238:GLU:OE2	5:S:238:GLU:HA	2.13	0.48
13:a:223:TYR:OH	13:a:253:ASN:N	2.47	0.48
5:E:62:LYS:NZ	5:E:74:ILE:O	2.39	0.48
5:E:121:GLN:NE2	6:F:130:ARG:O	2.46	0.48
13:a:238:THR:OG1	13:a:239:GLU:N	2.46	0.48
15:c:233:TYR:CG	15:c:253:ILE:HD11	2.49	0.47
14:N:20:THR:HG23	14:N:31:THR:HG21	1.96	0.47
3:Q:192:ILE:HD11	3:Q:208:LEU:HD13	1.96	0.47
15:c:84:SER:HA	15:c:88:MET:HG2	1.96	0.47
1:A:69:THR:OG1	1:A:70:LYS:N	2.48	0.47
13:M:238:THR:OG1	13:M:239:GLU:N	2.44	0.47
15:f:132:VAL:HG11	15:f:262:LEU:HD11	1.96	0.47
8:V:-26:ASN:OD1	8:V:-26:ASN:N	2.47	0.47
1:O:69:THR:OG1	1:O:70:LYS:N	2.47	0.47
6:T:98:ASN:ND2	17:h:68:ARG:NH1	2.63	0.47
15:c:201:ALA:O	15:c:202:CYS:SG	2.72	0.47
1:A:119:GLN:OE1	17:e:91:ARG:NH2	2.45	0.47
1:A:195:LEU:O	1:A:199:PHE:N	2.48	0.47
4:D:59:MET:SD	4:D:64:ILE:HD11	2.55	0.47
13:M:54:THR:OG1	13:M:227:ARG:O	2.33	0.47
13:a:86:ARG:NH1	13:a:98:ALA:O	2.48	0.47
1:A:44:VAL:HG11	1:A:137:CYS:HB2	1.96	0.47
4:R:175:GLU:OE1	4:R:175:GLU:N	2.38	0.47
12:Z:173:LEU:HD21	12:Z:210:ALA:HB2	1.97	0.47
3:C:91:CYS:SG	3:C:107:ILE:HD13	2.55	0.47
3:C:105:GLU:OE2	3:C:145:TYR:CZ	2.68	0.47
1:O:49:GLU:O	1:O:64:LYS:NZ	2.48	0.47
3:Q:80:ALA:HA	3:Q:129:ILE:HD13	1.96	0.47
11:Y:139:MET:O	11:Y:143:TYR:N	2.48	0.47
7:U:153:LYS:HB2	7:U:166:THR:HG21	1.96	0.46
17:e:89:VAL:HG13	17:e:89:VAL:O	2.13	0.46
11:K:139:MET:O	11:K:143:TYR:N	2.48	0.46
6:F:54:VAL:HG23	6:F:54:VAL:O	2.14	0.46
6:T:54:VAL:HG23	6:T:54:VAL:O	2.15	0.46
12:Z:56:ARG:NE	12:Z:219:ASP:OD1	2.37	0.46
8:H:1:THR:HG23	8:H:33:LYS:CD	2.46	0.46
8:H:166:ASP:OD1	8:H:168:GLY:N	2.45	0.46
6:T:153:ASP:OD1	6:T:154:PRO:N	2.48	0.46
3:Q:188:ILE:O	3:Q:192:ILE:HG12	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:80:ASP:OD2	5:S:126:ARG:NH1	2.42	0.46
6:T:127:SER:O	7:U:3:ARG:NH2	2.49	0.46
13:a:108:LEU:HD21	13:a:151:LEU:HD13	1.97	0.46
15:f:88:MET:N	15:f:88:MET:HE2	2.31	0.46
1:O:52:GLN:OE1	1:O:52:GLN:HA	2.16	0.46
15:c:88:MET:HE2	15:c:88:MET:N	2.30	0.46
3:C:166:LYS:O	3:C:170:GLU:HG3	2.16	0.46
6:T:98:ASN:CG	17:h:68:ARG:NH1	2.74	0.46
16:g:214:ILE:HD11	16:g:254:LEU:HD11	1.98	0.46
4:R:68:VAL:HG21	4:R:89:ILE:HG13	1.97	0.46
8:V:1:THR:HG23	8:V:33:LYS:HD3	1.97	0.46
15:f:210:ASN:O	15:f:211:ILE:HD13	2.15	0.46
12:Z:153:ASP:OD1	12:Z:157:SER:N	2.49	0.46
12:Z:169:ALA:O	12:Z:173:LEU:HG	2.16	0.46
13:a:211:ARG:NH1	13:a:245:GLU:OE2	2.42	0.46
9:W:30:ILE:HD12	9:W:30:ILE:H	1.79	0.45
4:D:85:ALA:O	4:D:89:ILE:HG12	2.16	0.45
4:D:85:ALA:HB2	4:D:139:VAL:HG21	1.97	0.45
4:D:175:GLU:OE1	4:D:175:GLU:N	2.39	0.45
13:M:108:LEU:HD21	13:M:151:LEU:HD13	1.98	0.45
3:Q:39:ASP:OD1	3:Q:213:ARG:NE	2.50	0.45
5:S:26:MET:CE	16:g:263:LEU:HD21	2.46	0.45
4:D:203:LYS:HB2	4:D:210:LEU:HD22	1.98	0.45
5:S:121:GLN:NE2	6:T:130:ARG:O	2.50	0.45
7:G:174:GLU:OE2	7:G:174:GLU:N	2.31	0.45
12:L:169:ALA:O	12:L:173:LEU:HG	2.16	0.45
7:U:115:CYS:O	7:U:115:CYS:SG	2.74	0.45
8:V:172:ASN:OD1	8:V:191:VAL:HG13	2.16	0.45
15:f:84:SER:HA	15:f:88:MET:HG2	1.98	0.45
11:Y:97:MET:HE3	11:Y:99:THR:HG22	1.99	0.45
13:M:104:ASP:O	13:M:108:LEU:HD23	2.16	0.45
3:Q:119:THR:HG22	3:Q:126:PRO:HB3	1.99	0.45
16:d:11:ASP:OD1	16:d:11:ASP:O	2.35	0.45
1:A:74:LEU:HD21	1:A:87:VAL:HG22	1.99	0.45
2:B:165:GLY:O	2:B:168:SER:OG	2.28	0.45
1:A:119:GLN:HA	1:A:122:THR:HG22	1.99	0.45
1:A:181:ASP:OD1	1:A:181:ASP:C	2.59	0.45
5:E:81:ALA:HB2	5:E:130:VAL:HG21	1.97	0.45
6:T:153:ASP:OD1	6:T:153:ASP:C	2.59	0.45
3:C:105:GLU:OE2	3:C:105:GLU:HA	2.17	0.45
7:G:138:MET:O	7:G:154:CYS:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:154:GLU:H	10:J:154:GLU:CD	2.24	0.45
12:L:229:GLU:OE1	12:L:229:GLU:N	2.45	0.45
13:M:111:VAL:O	13:M:115:MET:HG3	2.17	0.45
6:T:98:ASN:OD1	6:T:98:ASN:C	2.60	0.45
7:U:8:GLY:O	7:U:15:ILE:HD11	2.17	0.45
3:C:116:GLN:O	3:C:119:THR:OG1	2.25	0.45
8:H:1:THR:HG23	8:H:33:LYS:HD3	1.99	0.45
7:U:88:ARG:O	7:U:92:GLN:HG2	2.17	0.45
16:d:154:LYS:O	16:d:157:SER:OG	2.28	0.45
16:g:179:ILE:O	16:g:179:ILE:CG2	2.65	0.45
2:P:8:ARG:HB2	2:P:11:ILE:HD11	1.98	0.44
2:P:90:LEU:HG	2:P:114:LEU:HD13	1.99	0.44
15:f:244:LEU:O	15:f:247:VAL:HG22	2.16	0.44
3:Q:120:GLN:OE1	3:Q:120:GLN:HA	2.16	0.44
9:W:138:VAL:HG22	9:W:147:TYR:CE1	2.52	0.44
3:C:175:ASN:OD1	3:C:175:ASN:N	2.50	0.44
12:L:153:ASP:OD1	12:L:157:SER:N	2.51	0.44
10:X:161:ARG:O	10:X:165:GLU:HG3	2.17	0.44
15:c:285:ASN:OD1	15:c:285:ASN:N	2.50	0.44
8:V:-2:THR:O	8:V:33:LYS:NZ	2.51	0.44
16:g:181:ILE:HD12	16:g:187:THR:HG23	1.99	0.44
3:Q:175:ASN:N	3:Q:175:ASN:OD1	2.50	0.44
5:S:88:MET:HE3	5:S:112:ILE:HD11	2.00	0.44
8:V:-2:THR:O	8:V:-1:GLY:C	2.59	0.44
13:a:111:VAL:O	13:a:115:MET:HG3	2.18	0.44
6:F:18:ASP:OD1	6:F:18:ASP:N	2.37	0.44
6:T:24:VAL:O	6:T:28:MET:HG3	2.18	0.44
14:b:20:THR:HG23	14:b:31:THR:HG21	1.99	0.44
7:U:53:GLN:NE2	7:U:206:LEU:HD11	2.32	0.44
2:B:183:GLU:OE1	2:B:183:GLU:N	2.48	0.44
6:F:34:SER:O	6:F:168:LYS:N	2.51	0.44
7:G:80:MET:SD	7:G:138:MET:HG3	2.58	0.44
1:O:119:GLN:HA	1:O:122:THR:HG22	1.99	0.44
8:V:166:ASP:OD1	8:V:168:GLY:N	2.48	0.44
16:d:20:PRO:HG2	16:d:90:LEU:H	1.83	0.44
12:L:173:LEU:HD22	12:L:206:VAL:HB	2.00	0.44
10:J:11:ASP:N	10:J:11:ASP:OD1	2.50	0.43
8:V:59:ILE:HD11	8:V:86:MET:HE2	2.00	0.43
15:f:144:ASP:O	15:f:148:GLN:NE2	2.42	0.43
7:G:138:MET:HB3	7:G:154:CYS:HB3	1.99	0.43
2:P:46:ALA:HB1	2:P:197:LEU:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:69:ASN:OD1	6:T:225:HIS:ND1	2.49	0.43
1:O:74:LEU:HD21	1:O:87:VAL:HG22	2.01	0.43
3:Q:192:ILE:O	3:Q:196:LEU:HD23	2.18	0.43
2:B:112:THR:HG23	3:C:81:ARG:HD2	2.00	0.43
4:D:68:VAL:HG21	4:D:89:ILE:HG13	2.00	0.43
5:S:26:MET:O	5:S:29:VAL:HG12	2.18	0.43
7:U:115:CYS:SG	7:U:160:TYR:HB2	2.58	0.43
10:X:11:ASP:N	10:X:11:ASP:OD1	2.51	0.43
17:h:75:ALA:HB3	17:h:76:PRO:HD3	2.01	0.43
3:C:119:THR:HG22	3:C:126:PRO:HB3	2.01	0.43
8:H:55:THR:HG23	8:H:86:MET:HE1	1.99	0.43
13:M:223:TYR:OH	13:M:253:ASN:N	2.52	0.43
1:O:195:LEU:O	1:O:199:PHE:N	2.51	0.43
13:a:83:ASN:OD1	13:a:231:ARG:NH2	2.49	0.43
5:E:26:MET:O	5:E:29:VAL:HG12	2.18	0.43
7:G:8:GLY:O	7:G:15:ILE:HD11	2.18	0.43
7:G:206:LEU:HB3	7:G:208:ILE:HG22	2.00	0.43
8:H:104:ASP:OD1	8:H:106:THR:OG1	2.31	0.43
8:H:153:ASN:O	8:H:157:GLU:HG3	2.18	0.43
14:N:166:ARG:NE	14:b:140:ASP:OD2	2.48	0.43
12:Z:173:LEU:HD22	12:Z:206:VAL:HB	2.01	0.43
13:a:104:ASP:O	13:a:108:LEU:HD23	2.18	0.43
1:A:49:GLU:O	1:A:64:LYS:NZ	2.52	0.43
10:X:13:VAL:HG13	10:X:13:VAL:O	2.18	0.43
15:f:93:TRP:HZ2	15:f:262:LEU:HD12	1.84	0.43
2:B:123:GLN:NE2	3:C:125:ARG:O	2.52	0.43
14:N:144:ARG:NH2	14:N:151:GLU:OE2	2.49	0.43
10:X:164:LEU:O	10:X:168:GLN:HG2	2.18	0.43
15:f:82:PHE:HB2	15:f:247:VAL:HG12	2.01	0.43
6:F:153:ASP:OD1	6:F:154:PRO:HD2	2.18	0.43
12:L:173:LEU:HD21	12:L:210:ALA:HB2	2.01	0.43
4:R:161:THR:HG22	5:S:78:THR:HB	2.00	0.43
7:G:201:CYS:O	7:G:205:VAL:HG22	2.18	0.42
2:P:112:THR:HG23	3:Q:81:ARG:HD2	2.00	0.42
2:P:174:MET:SD	2:P:199:LYS:HG2	2.59	0.42
3:Q:39:ASP:OD1	3:Q:213:ARG:NH2	2.52	0.42
4:R:85:ALA:O	4:R:89:ILE:HG12	2.18	0.42
7:U:123:GLN:NE2	7:U:127:GLN:OE1	2.47	0.42
8:V:-27:ASP:OD1	8:V:-27:ASP:N	2.52	0.42
2:B:90:LEU:HG	2:B:114:LEU:HD13	2.01	0.42
14:b:62:GLN:OE1	14:b:86:MET:HE1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:161:THR:HG22	5:E:78:THR:HB	2.01	0.42
4:R:85:ALA:HB2	4:R:139:VAL:HG21	1.99	0.42
8:V:166:ASP:OD1	8:V:167:LEU:N	2.53	0.42
16:d:11:ASP:OD1	16:d:11:ASP:C	2.62	0.42
16:d:142:LEU:HD21	16:d:179:ILE:HG12	2.01	0.42
17:e:45:SER:N	17:e:49:GLU:OE1	2.53	0.42
3:C:98:VAL:O	3:C:99:GLU:HB3	2.19	0.42
12:L:96:ILE:HD11	12:L:120:LEU:HD13	2.01	0.42
2:P:174:MET:HE2	2:P:196:VAL:HG22	2.00	0.42
3:Q:98:VAL:O	3:Q:99:GLU:HB3	2.19	0.42
12:Z:142:ASP:OD1	12:Z:146:LYS:N	2.50	0.42
12:Z:176:LEU:HD23	12:Z:206:VAL:CG1	2.49	0.42
16:g:11:ASP:OD1	16:g:11:ASP:O	2.37	0.42
8:H:-2:THR:O	8:H:-1:GLY:C	2.61	0.42
8:V:1:THR:HG23	8:V:33:LYS:CD	2.50	0.42
17:h:67:LEU:O	17:h:71:GLN:N	2.51	0.42
1:A:177:ARG:HH11	1:A:177:ARG:HG2	1.84	0.42
7:G:122:SER:O	7:G:126:THR:HG23	2.20	0.42
9:I:34:MET:O	11:Y:166:ARG:NH1	2.52	0.42
2:P:82:ASP:OD1	2:P:82:ASP:N	2.53	0.42
10:X:92:LEU:HD12	10:X:97:PRO:HB3	2.01	0.42
17:e:48:LEU:O	17:e:52:GLU:HG2	2.19	0.42
5:E:22:ILE:O	5:E:26:MET:HG3	2.19	0.42
8:H:26:VAL:O	12:Z:213:ARG:NH1	2.53	0.42
10:J:164:LEU:O	10:J:168:GLN:HG2	2.19	0.42
14:N:59:VAL:HG11	14:N:83:PHE:CE2	2.55	0.42
14:N:205:ALA:HB3	13:a:83:ASN:HD22	1.84	0.42
6:T:18:ASP:OD1	6:T:18:ASP:N	2.45	0.42
16:d:229:GLN:OE1	16:d:229:GLN:HA	2.20	0.42
17:h:68:ARG:NE	17:h:73:LEU:HG	2.35	0.42
16:d:107:LEU:HD22	16:d:199:ILE:HD11	2.01	0.42
16:d:120:VAL:HB	16:d:203:VAL:HG22	2.02	0.42
10:J:27:GLN:OE1	10:J:174:ASN:OD1	2.37	0.42
1:O:184:LEU:O	1:O:185:GLU:OE1	2.38	0.42
2:P:8:ARG:O	2:P:9:THR:OG1	2.33	0.42
3:Q:91:CYS:SG	3:Q:107:ILE:HD13	2.60	0.42
13:a:96:LEU:HD12	13:a:157:ILE:HG12	2.01	0.42
1:A:184:LEU:C	1:A:186:ASP:H	2.28	0.41
3:C:171:PHE:CD1	3:C:171:PHE:C	2.98	0.41
4:D:28:ILE:HD13	4:D:28:ILE:N	2.35	0.41
5:E:80:ASP:OD2	5:E:126:ARG:NH1	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:34:SER:O	6:T:168:LYS:N	2.53	0.41
10:X:49:GLU:O	10:X:50:ALA:C	2.63	0.41
17:h:89:VAL:O	17:h:89:VAL:CG1	2.67	0.41
1:A:127:VAL:HG13	17:e:89:VAL:HG11	2.02	0.41
11:K:37:ILE:HG23	11:K:60:ALA:HA	2.02	0.41
12:Z:76:ASP:OD1	12:Z:76:ASP:O	2.38	0.41
16:g:148:GLN:O	16:g:151:VAL:HG22	2.20	0.41
10:J:13:VAL:O	10:J:13:VAL:HG13	2.20	0.41
12:L:212:GLU:OE2	12:L:239:ARG:NH1	2.46	0.41
2:P:35:LEU:HD23	2:P:35:LEU:C	2.45	0.41
3:Q:4:ASP:OD2	3:Q:21:TYR:OH	2.38	0.41
15:c:70:LYS:HG2	15:c:72:ILE:HD11	2.02	0.41
8:H:112:SER:O	8:H:112:SER:OG	2.38	0.41
10:J:49:GLU:O	10:J:50:ALA:C	2.63	0.41
14:N:140:ASP:OD2	14:b:166:ARG:NE	2.50	0.41
4:R:70:ILE:HD11	4:R:89:ILE:HD12	2.02	0.41
15:f:266:ILE:O	15:f:266:ILE:HG23	2.21	0.41
3:C:7:ILE:H	3:C:7:ILE:HD12	1.85	0.41
7:G:78:CYS:HB3	7:G:140:LEU:HD23	2.01	0.41
10:J:7:ILE:O	10:J:14:LEU:N	2.49	0.41
10:J:174:ASN:HB3	10:X:174:ASN:HB3	2.01	0.41
14:N:203:PRO:N	14:N:204:PRO:HD2	2.35	0.41
8:V:-2:THR:OG1	8:V:1:THR:OG1	2.37	0.41
17:e:68:ARG:HD2	17:e:68:ARG:O	2.19	0.41
7:G:31:ALA:HB1	7:G:83:MET:HE3	2.03	0.41
12:Z:52:ALA:HB1	12:Z:221:LEU:HD11	2.03	0.41
16:d:37:SER:O	16:d:246:LYS:NZ	2.49	0.41
17:e:89:VAL:O	17:e:89:VAL:CG1	2.69	0.41
16:g:120:VAL:HB	16:g:203:VAL:HG22	2.03	0.41
1:A:114:VAL:O	1:A:118:MET:HG2	2.20	0.41
14:N:2:THR:HG21	14:N:163:ALA:CB	2.50	0.41
1:O:108:ALA:O	1:O:112:GLN:HG3	2.20	0.41
2:P:206:LEU:N	2:P:206:LEU:HD23	2.35	0.41
5:S:60:GLN:OE1	15:f:288:THR:HB	2.20	0.41
11:Y:154:ASP:OD1	11:Y:154:ASP:N	2.53	0.41
1:A:228:ASP:OD1	1:A:228:ASP:N	2.54	0.41
5:E:11:THR:OG1	7:G:3:ARG:NH1	2.53	0.41
10:J:118:MET:HE2	10:J:124:LEU:HD13	2.03	0.41
2:P:231:LYS:O	2:P:235:GLN:HG2	2.21	0.41
9:W:159:ASP:OD1	9:W:160:PRO:HD2	2.21	0.41
9:W:161:ASP:OD1	9:W:161:ASP:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:c:19:GLU:OE1	15:c:19:GLU:N	2.52	0.41
15:c:244:LEU:O	15:c:247:VAL:HG22	2.21	0.41
6:F:69:ASN:OD1	6:F:225:HIS:ND1	2.50	0.41
12:Z:188:ASN:O	12:Z:188:ASN:ND2	2.54	0.41
16:d:230:ILE:C	16:d:231:LEU:HD12	2.46	0.41
2:B:34:CYS:SG	2:B:75:SER:OG	2.55	0.40
8:V:163:ILE:HG23	8:V:170:GLY:HA2	2.03	0.40
14:b:1:THR:HG22	14:b:2:THR:N	2.36	0.40
14:b:203:PRO:N	14:b:204:PRO:HD2	2.35	0.40
16:d:148:GLN:O	16:d:152:GLN:HG3	2.21	0.40
6:T:51:GLU:OE2	6:T:202:HIS:ND1	2.32	0.40
12:Z:207:PHE:CG	12:Z:221:LEU:HD13	2.56	0.40
1:A:40:ALA:N	1:A:43:GLY:O	2.53	0.40
4:R:225:ASN:OD1	4:R:225:ASN:C	2.63	0.40
5:S:209:ASN:OD1	5:S:209:ASN:N	2.53	0.40
6:T:188:ARG:O	6:T:191:VAL:HG12	2.21	0.40
8:V:42:TYR:HB2	8:V:178:ILE:HD11	2.04	0.40
9:W:53:LEU:HB2	9:W:60:VAL:HG13	2.02	0.40
8:H:214:GLU:HA	8:H:214:GLU:OE1	2.21	0.40
1:O:40:ALA:N	1:O:43:GLY:O	2.54	0.40
15:c:165:THR:CG2	15:c:167:LEU:HD21	2.51	0.40
16:d:230:ILE:O	16:d:231:LEU:HD12	2.22	0.40
17:h:45:SER:N	17:h:49:GLU:OE1	2.52	0.40
3:C:96:LEU:HB2	10:J:62:LYS:HG3	2.03	0.40
3:C:117:ARG:HD3	17:e:140:LEU:HD13	2.03	0.40
7:G:88:ARG:O	7:G:92:GLN:HG2	2.21	0.40
2:P:122:THR:HG22	2:P:129:PRO:HB3	2.04	0.40
3:Q:155:ALA:HB3	4:R:63:SER:HB2	2.02	0.40
14:b:2:THR:HG21	14:b:163:ALA:CB	2.52	0.40
17:e:78:LYS:O	17:e:82:GLU:HG3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	232/234 (99%)	229 (99%)	3 (1%)	0	100	100
1	O	232/234 (99%)	225 (97%)	7 (3%)	0	100	100
2	B	249/261 (95%)	243 (98%)	6 (2%)	0	100	100
2	P	249/261 (95%)	244 (98%)	5 (2%)	0	100	100
3	C	236/248 (95%)	233 (99%)	3 (1%)	0	100	100
3	Q	236/248 (95%)	233 (99%)	3 (1%)	0	100	100
4	D	239/241 (99%)	235 (98%)	4 (2%)	0	100	100
4	R	239/241 (99%)	235 (98%)	4 (2%)	0	100	100
5	E	237/263 (90%)	233 (98%)	4 (2%)	0	100	100
5	S	237/263 (90%)	232 (98%)	5 (2%)	0	100	100
6	F	241/255 (94%)	239 (99%)	2 (1%)	0	100	100
6	T	241/255 (94%)	238 (99%)	3 (1%)	0	100	100
7	G	243/246 (99%)	241 (99%)	2 (1%)	0	100	100
7	U	243/246 (99%)	241 (99%)	2 (1%)	0	100	100
8	H	255/277 (92%)	244 (96%)	10 (4%)	1 (0%)	30	60
8	V	255/277 (92%)	243 (95%)	11 (4%)	1 (0%)	30	60
9	I	203/205 (99%)	195 (96%)	8 (4%)	0	100	100
9	W	203/205 (99%)	194 (96%)	9 (4%)	0	100	100
10	J	195/201 (97%)	190 (97%)	4 (2%)	1 (0%)	24	55
10	X	195/201 (97%)	190 (97%)	4 (2%)	1 (0%)	24	55
11	K	198/263 (75%)	196 (99%)	2 (1%)	0	100	100
11	Y	198/263 (75%)	196 (99%)	2 (1%)	0	100	100
12	L	211/241 (88%)	208 (99%)	3 (1%)	0	100	100
12	Z	211/241 (88%)	209 (99%)	2 (1%)	0	100	100
13	M	213/264 (81%)	207 (97%)	6 (3%)	0	100	100
13	a	213/264 (81%)	207 (97%)	6 (3%)	0	100	100
14	N	213/239 (89%)	210 (99%)	3 (1%)	0	100	100
14	b	213/239 (89%)	210 (99%)	3 (1%)	0	100	100
15	c	267/288 (93%)	256 (96%)	11 (4%)	0	100	100
15	f	267/288 (93%)	256 (96%)	11 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
16	d	243/264 (92%)	233 (96%)	10 (4%)	0	100	100
16	g	243/264 (92%)	235 (97%)	8 (3%)	0	100	100
17	e	101/141 (72%)	99 (98%)	2 (2%)	0	100	100
17	h	101/141 (72%)	98 (97%)	3 (3%)	0	100	100
All	All	7552/8262 (91%)	7377 (98%)	171 (2%)	4 (0%)	49	77

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
8	H	-40	VAL
8	V	-40	VAL
10	X	50	ALA
10	J	50	ALA

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	176/191 (92%)	176 (100%)	0	100	100
1	O	176/191 (92%)	176 (100%)	0	100	100
2	B	192/221 (87%)	191 (100%)	1 (0%)	81	93
2	P	192/221 (87%)	191 (100%)	1 (0%)	81	93
3	C	161/211 (76%)	161 (100%)	0	100	100
3	Q	161/211 (76%)	161 (100%)	0	100	100
4	D	184/203 (91%)	183 (100%)	1 (0%)	81	93
4	R	184/203 (91%)	184 (100%)	0	100	100
5	E	186/224 (83%)	186 (100%)	0	100	100
5	S	186/224 (83%)	186 (100%)	0	100	100
6	F	182/212 (86%)	181 (100%)	1 (0%)	81	93
6	T	182/212 (86%)	180 (99%)	2 (1%)	65	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	G	184/210 (88%)	184 (100%)	0	100	100
7	U	184/210 (88%)	184 (100%)	0	100	100
8	H	200/228 (88%)	200 (100%)	0	100	100
8	V	200/228 (88%)	200 (100%)	0	100	100
9	I	167/174 (96%)	167 (100%)	0	100	100
9	W	167/174 (96%)	167 (100%)	0	100	100
10	J	163/171 (95%)	163 (100%)	0	100	100
10	X	163/171 (95%)	163 (100%)	0	100	100
11	K	149/202 (74%)	149 (100%)	0	100	100
11	Y	149/202 (74%)	149 (100%)	0	100	100
12	L	171/199 (86%)	171 (100%)	0	100	100
12	Z	171/199 (86%)	171 (100%)	0	100	100
13	M	165/215 (77%)	165 (100%)	0	100	100
13	a	165/215 (77%)	165 (100%)	0	100	100
14	N	164/181 (91%)	164 (100%)	0	100	100
14	b	164/181 (91%)	164 (100%)	0	100	100
15	c	220/262 (84%)	220 (100%)	0	100	100
15	f	220/262 (84%)	220 (100%)	0	100	100
16	d	209/237 (88%)	209 (100%)	0	100	100
16	g	209/237 (88%)	209 (100%)	0	100	100
17	e	92/128 (72%)	92 (100%)	0	100	100
17	h	92/128 (72%)	92 (100%)	0	100	100
All	All	5930/6938 (86%)	5924 (100%)	6 (0%)	87	97

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	76	VAL
4	D	89	ILE
6	F	158	SER
2	P	76	VAL
6	T	82	LEU
6	T	102	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	GLN
1	A	102	GLN
1	A	148	GLN
2	B	51	ASN
2	B	102	GLN
2	B	119	GLN
3	C	116	GLN
4	D	155	HIS
5	E	90	GLN
7	G	12	HIS
9	I	173	ASN
10	J	24	ASN
10	J	101	ASN
10	J	174	ASN
11	K	29	GLN
12	L	179	ASN
12	L	180	GLN
1	O	96	GLN
1	O	102	GLN
1	O	148	GLN
2	P	30	HIS
2	P	51	ASN
2	P	102	GLN
2	P	155	ASN
3	Q	116	GLN
3	Q	205	ASN
4	R	13	ASN
4	R	152	GLN
4	R	204	GLN
4	R	227	HIS
5	S	185	ASN
6	T	121	HIS
7	U	12	HIS
7	U	172	GLN
9	W	173	ASN
10	X	27	GLN
10	X	101	ASN
10	X	174	ASN
12	Z	105	HIS
12	Z	179	ASN
12	Z	180	GLN

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Mol	Chain	Res	Type
12	Z	188	ASN
13	a	192	GLN
13	a	258	HIS
15	c	89	ASN
15	c	103	ASN
15	c	145	GLN
15	c	163	GLN
16	d	213	ASN
15	f	163	GLN
15	f	208	GLN
15	f	280	ASN
16	g	200	GLN
17	h	87	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

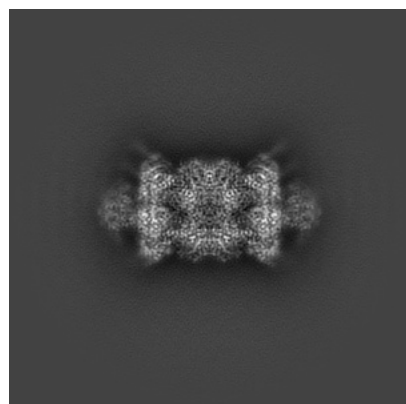
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41380. These allow visual inspection of the internal detail of the map and identification of artifacts.

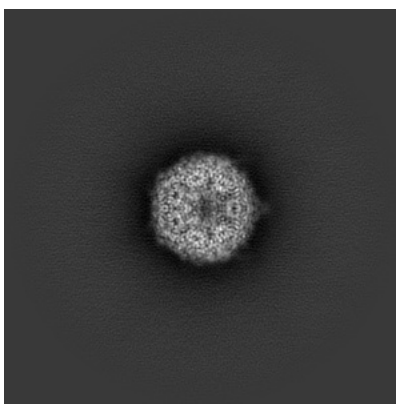
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

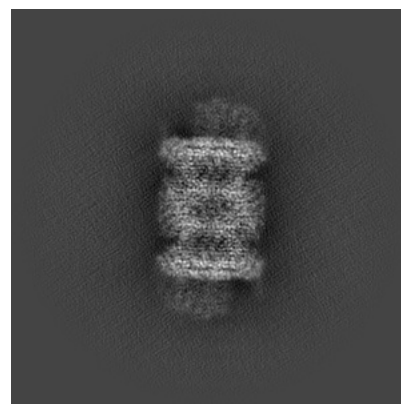
6.1.1 Primary map



X

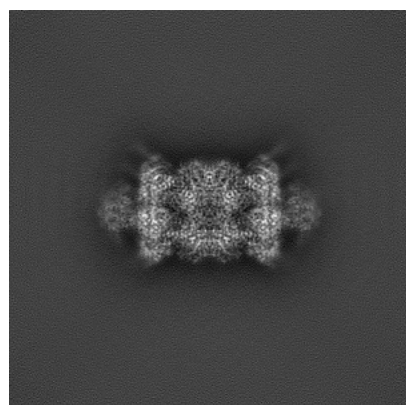


Y

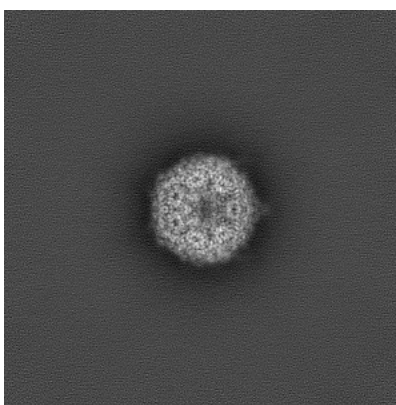


Z

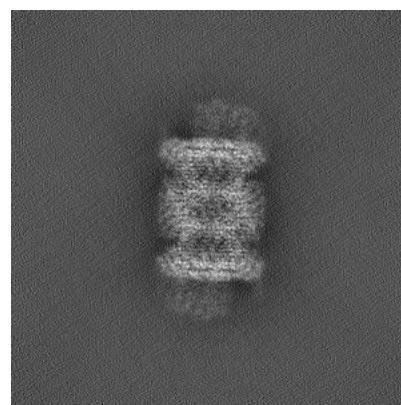
6.1.2 Raw map



X



Y

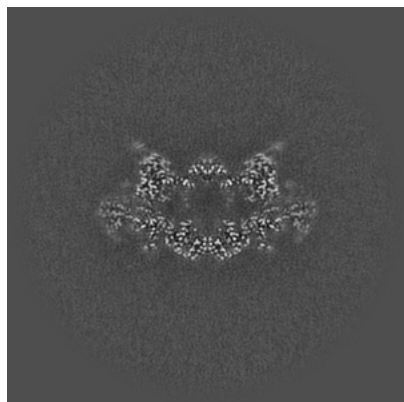


Z

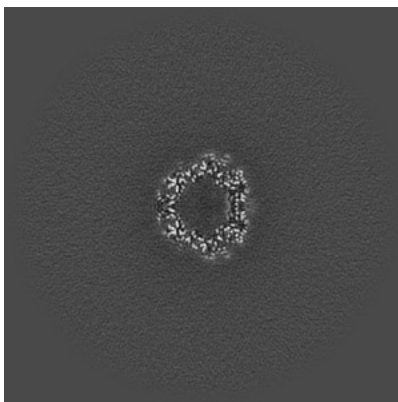
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

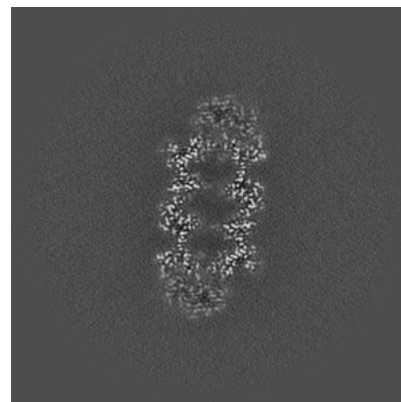
6.2.1 Primary map



X Index: 200

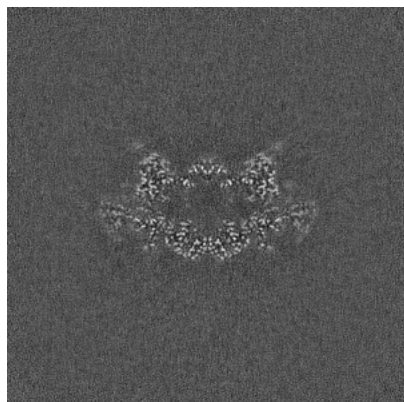


Y Index: 200

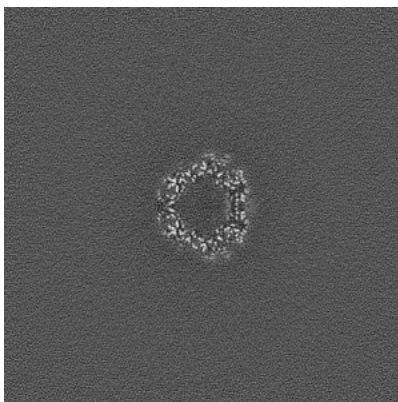


Z Index: 200

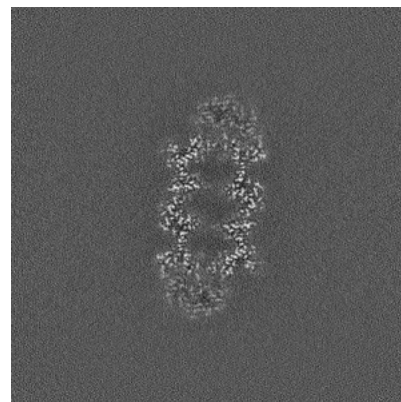
6.2.2 Raw map



X Index: 200



Y Index: 200

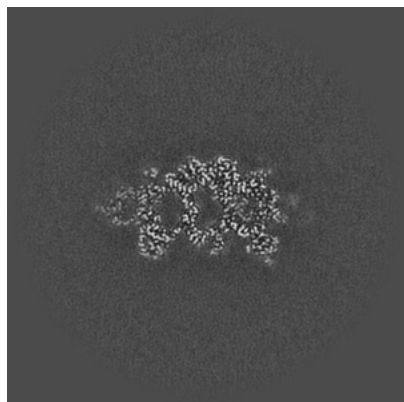


Z Index: 200

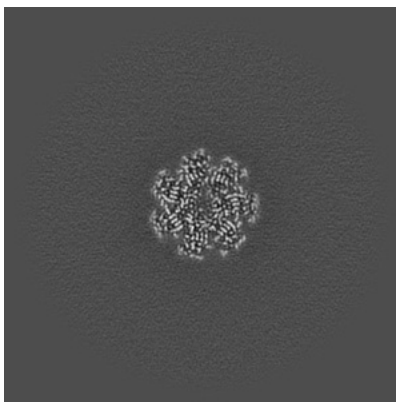
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

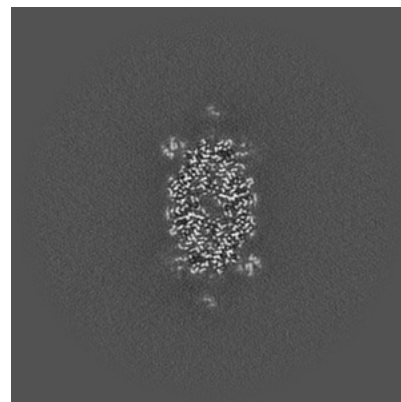
6.3.1 Primary map



X Index: 182

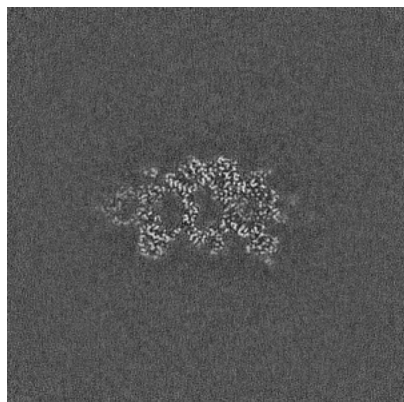


Y Index: 254

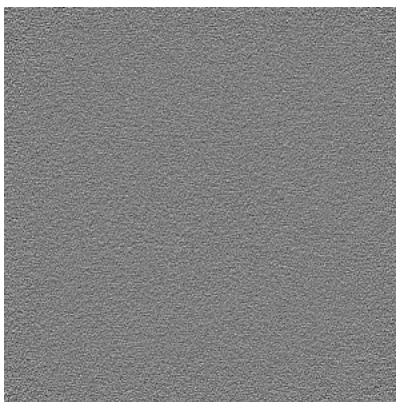


Z Index: 175

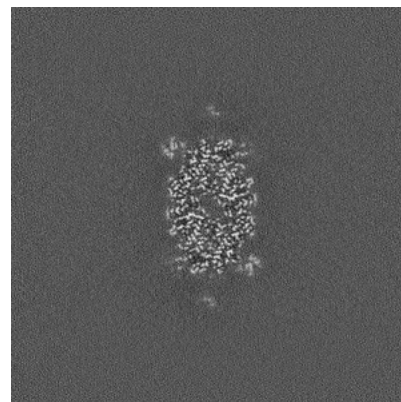
6.3.2 Raw map



X Index: 182



Y Index: 0

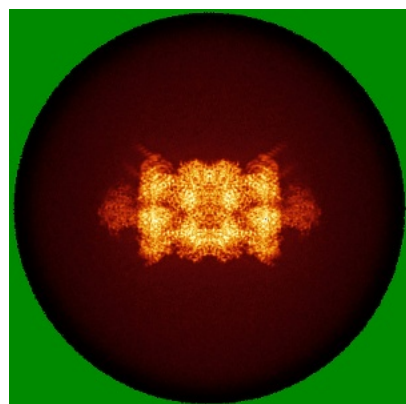


Z Index: 175

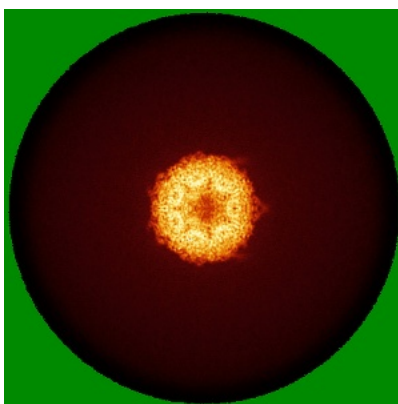
The images above show the largest variance slices of the map in three orthogonal directions.

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

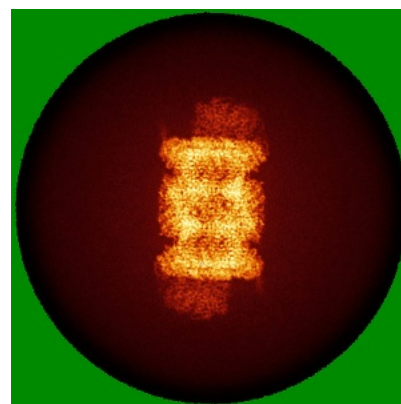
6.4.1 Primary map



X

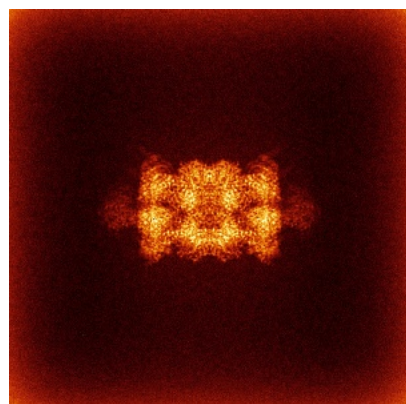


Y

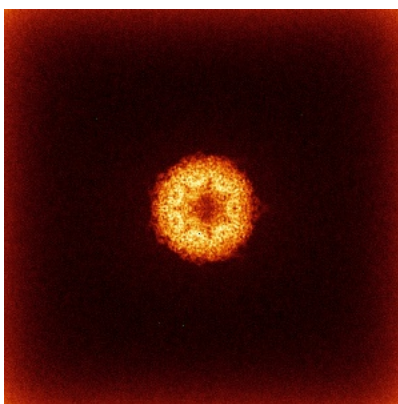


Z

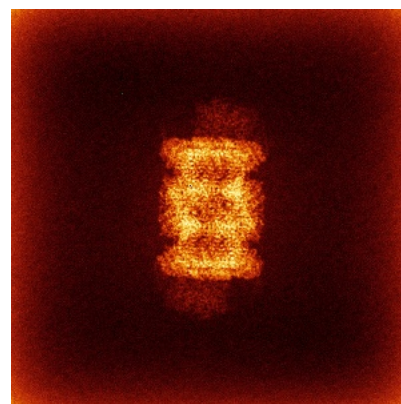
6.4.2 Raw map



X



Y

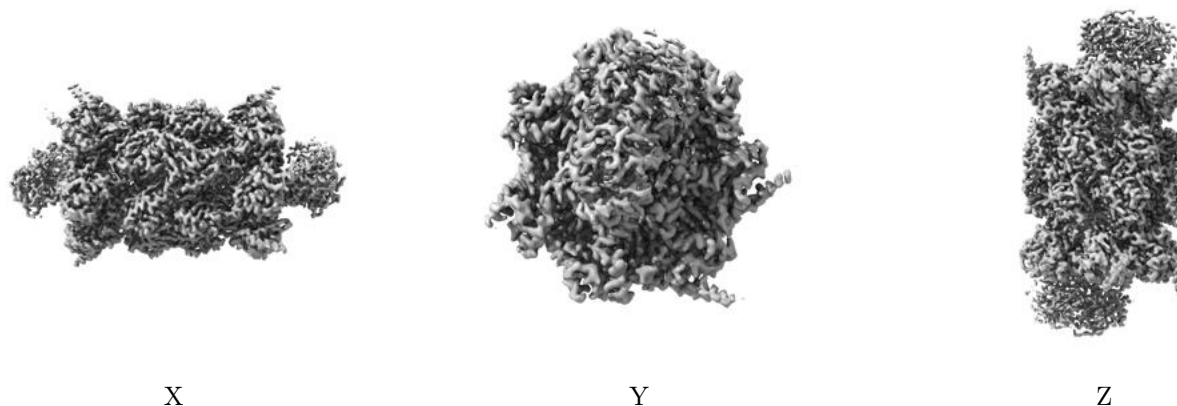


Z

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

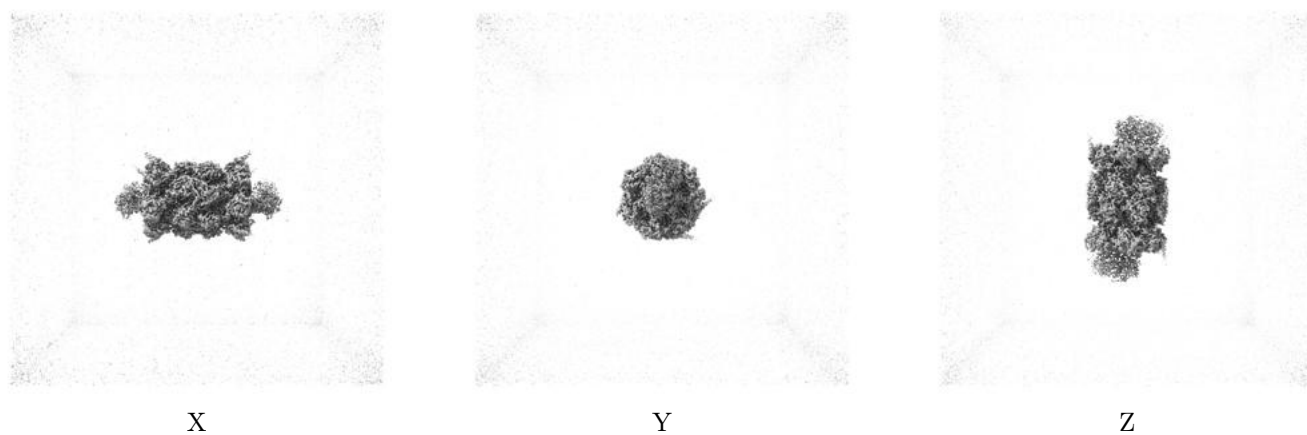
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 6.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

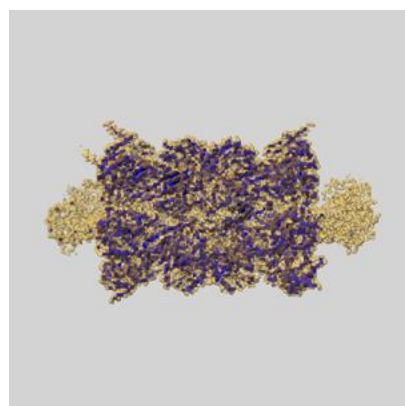
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

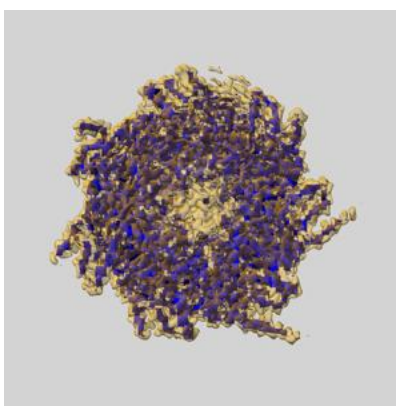
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

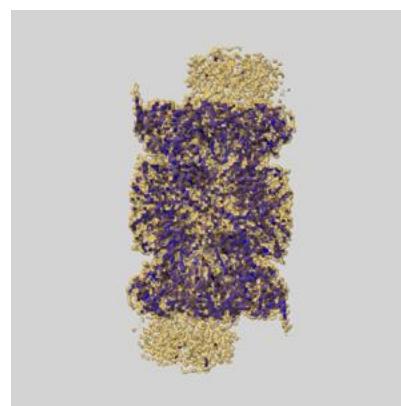
6.6.1 emd_41380_msk_1.map [i](#)



X



Y

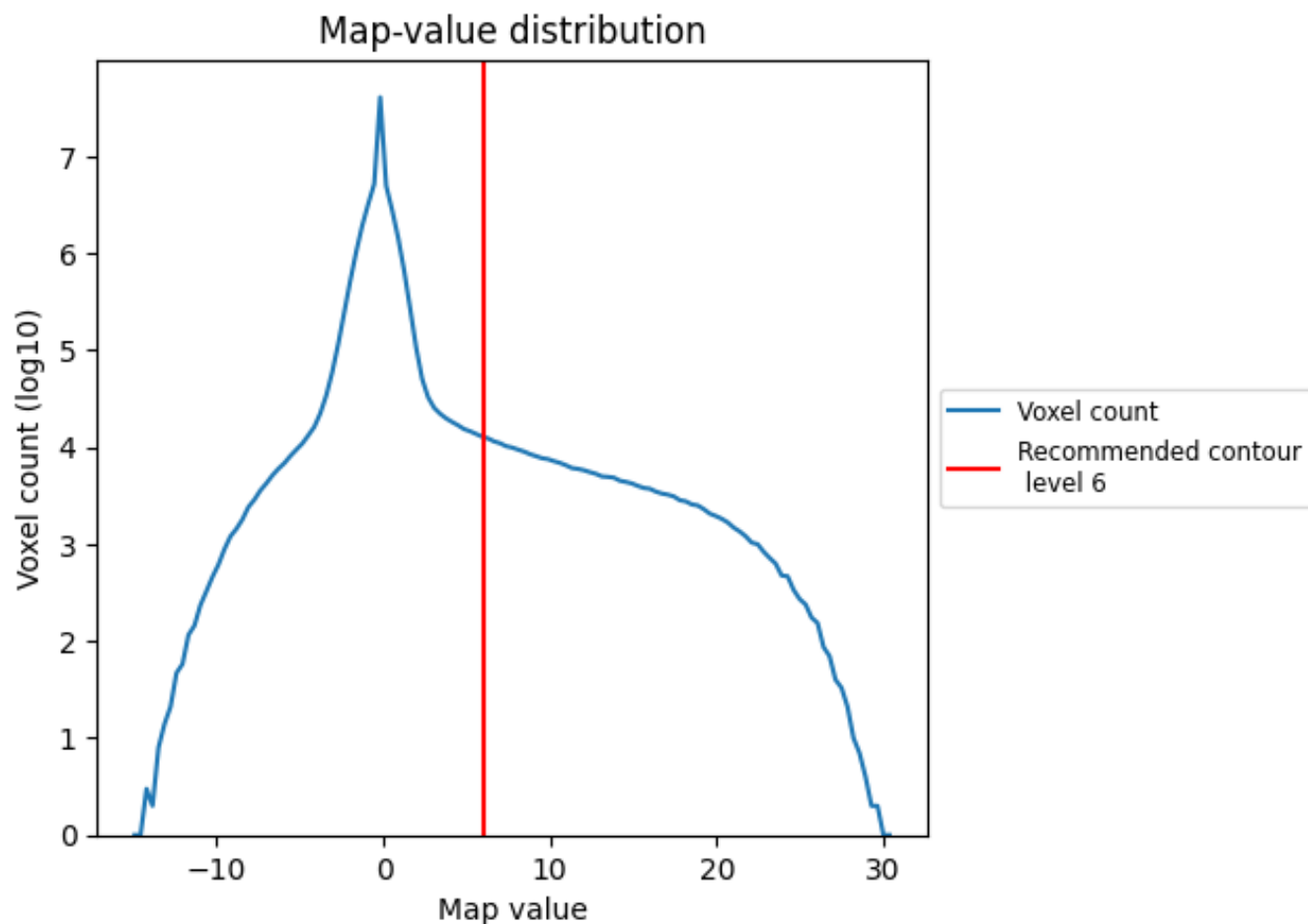


Z

7 Map analysis [i](#)

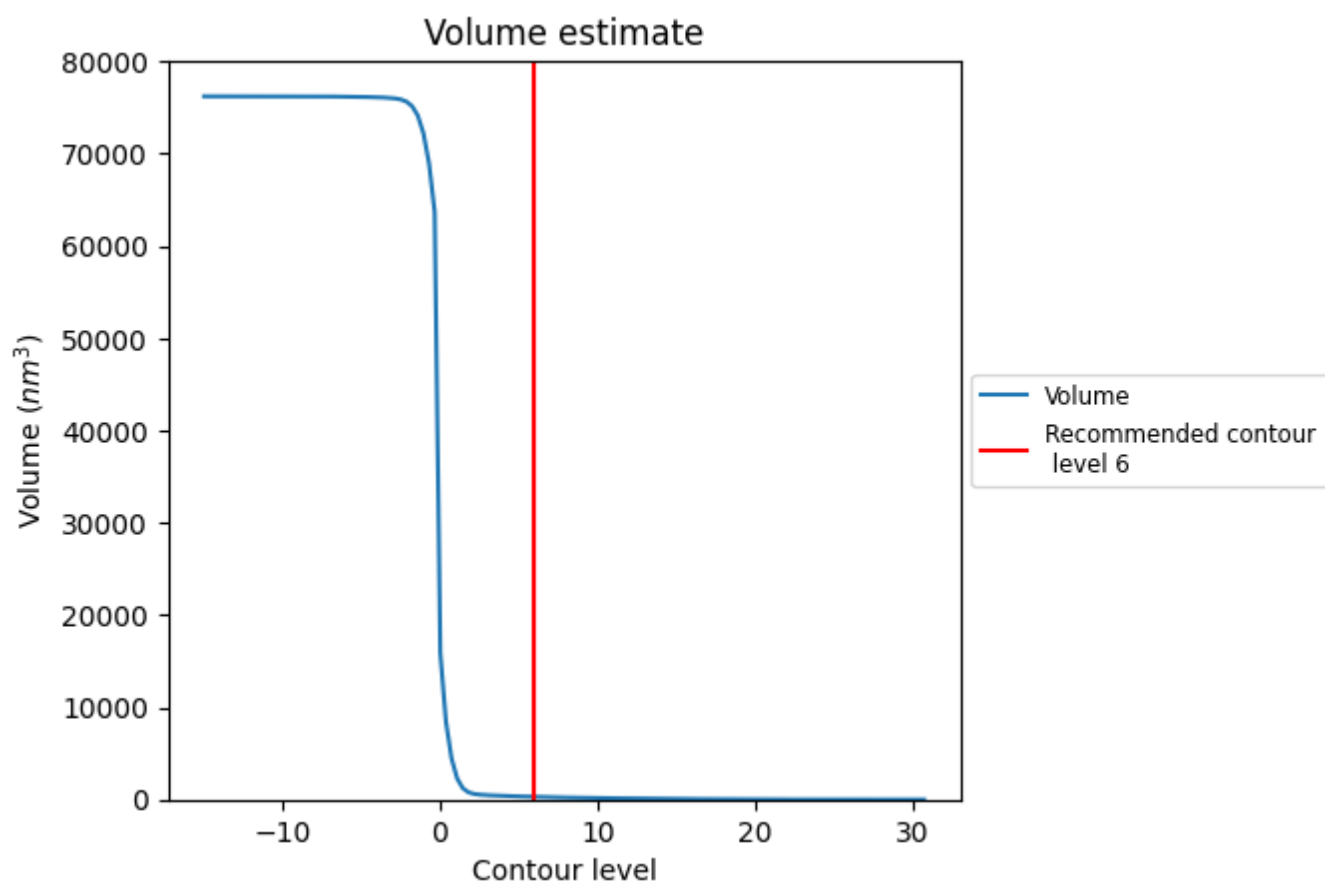
This section contains the results of statistical analysis of the map.

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.

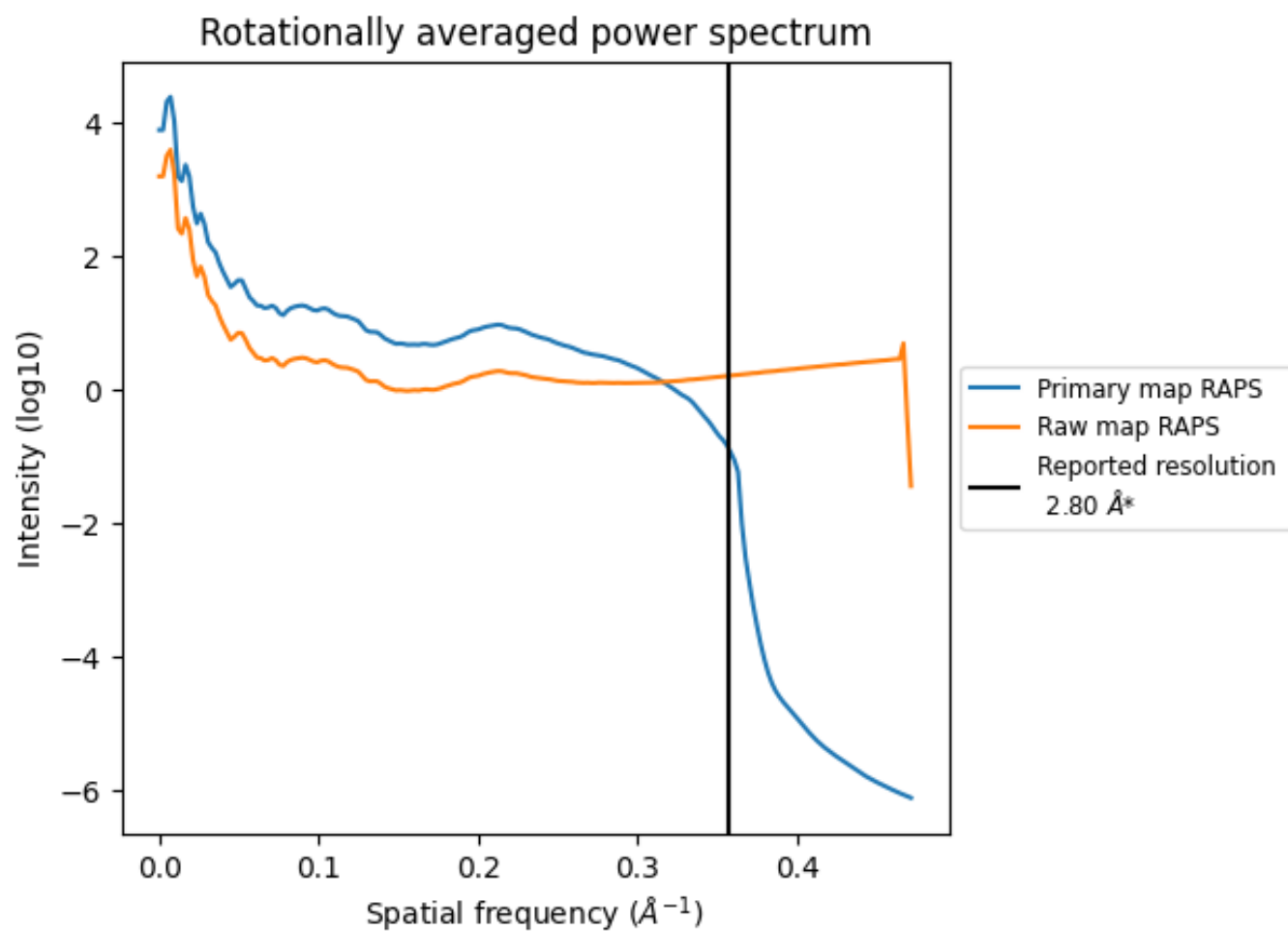
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 296 nm³; this corresponds to an approximate mass of 268 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 ⓘ

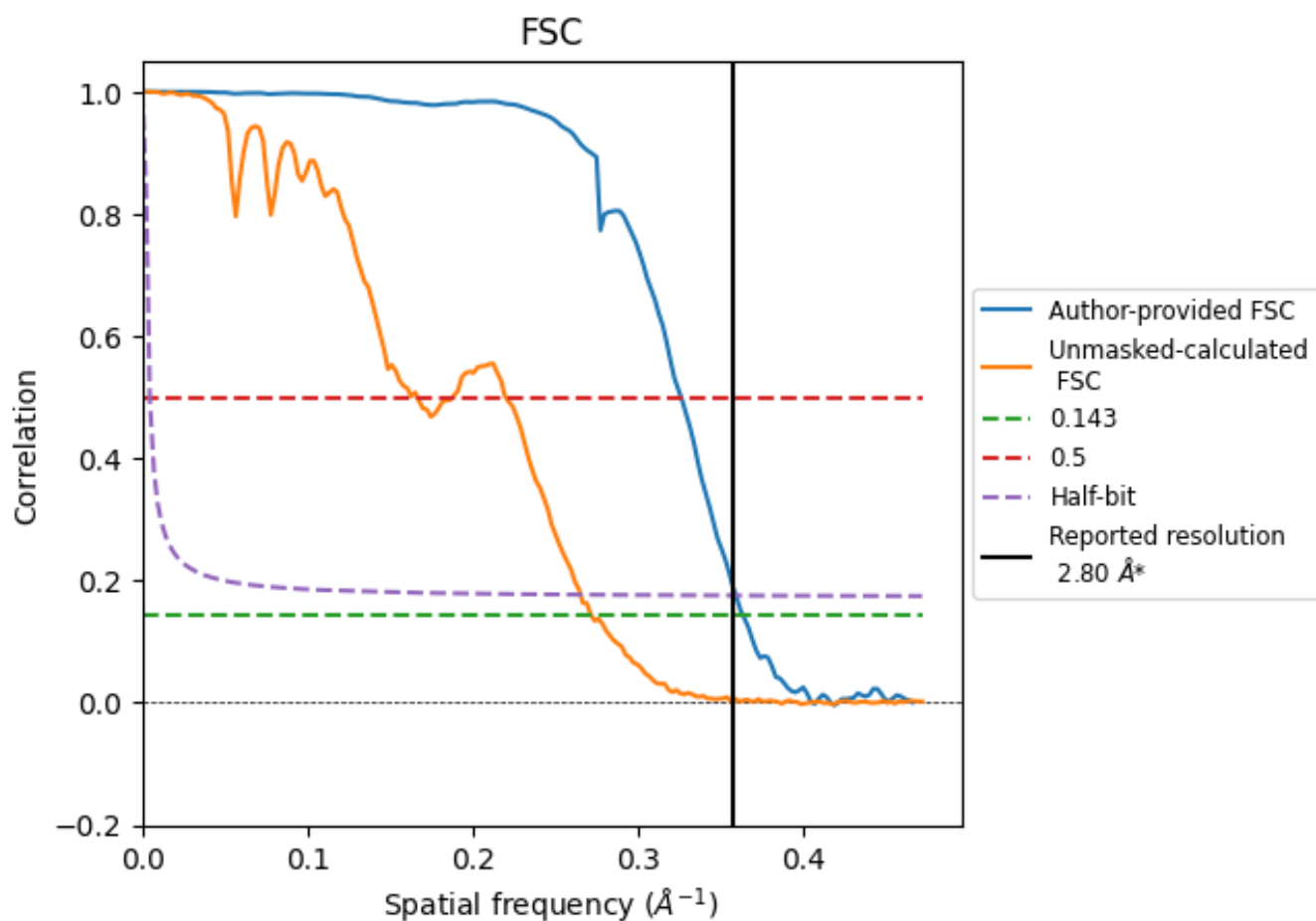


*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

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 0.357 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	2.75	3.07	2.79
Unmasked-calculated*	3.68	6.01	3.77

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.68 differs from the reported value 2.8 by more than 10 %

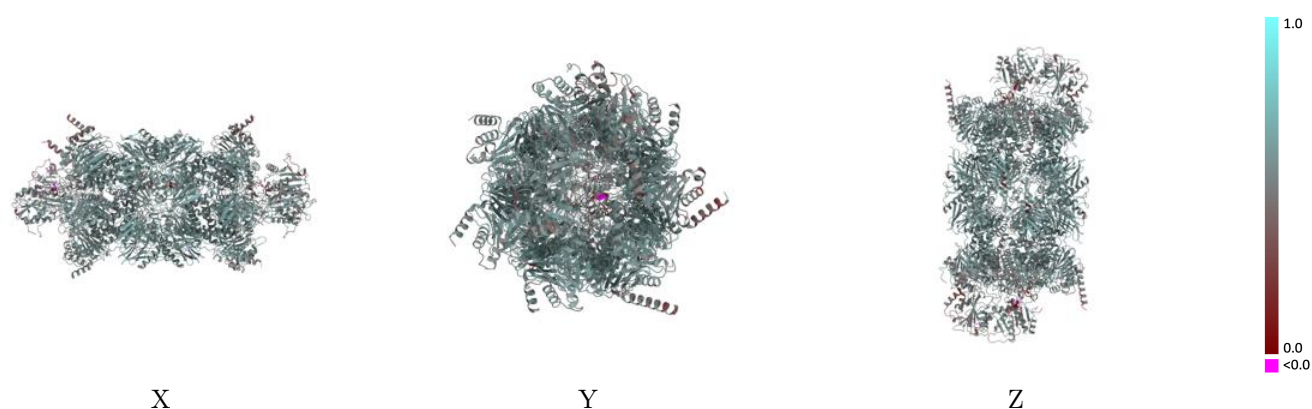
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-41380 and PDB model 8TM6. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)

This section was not generated.

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

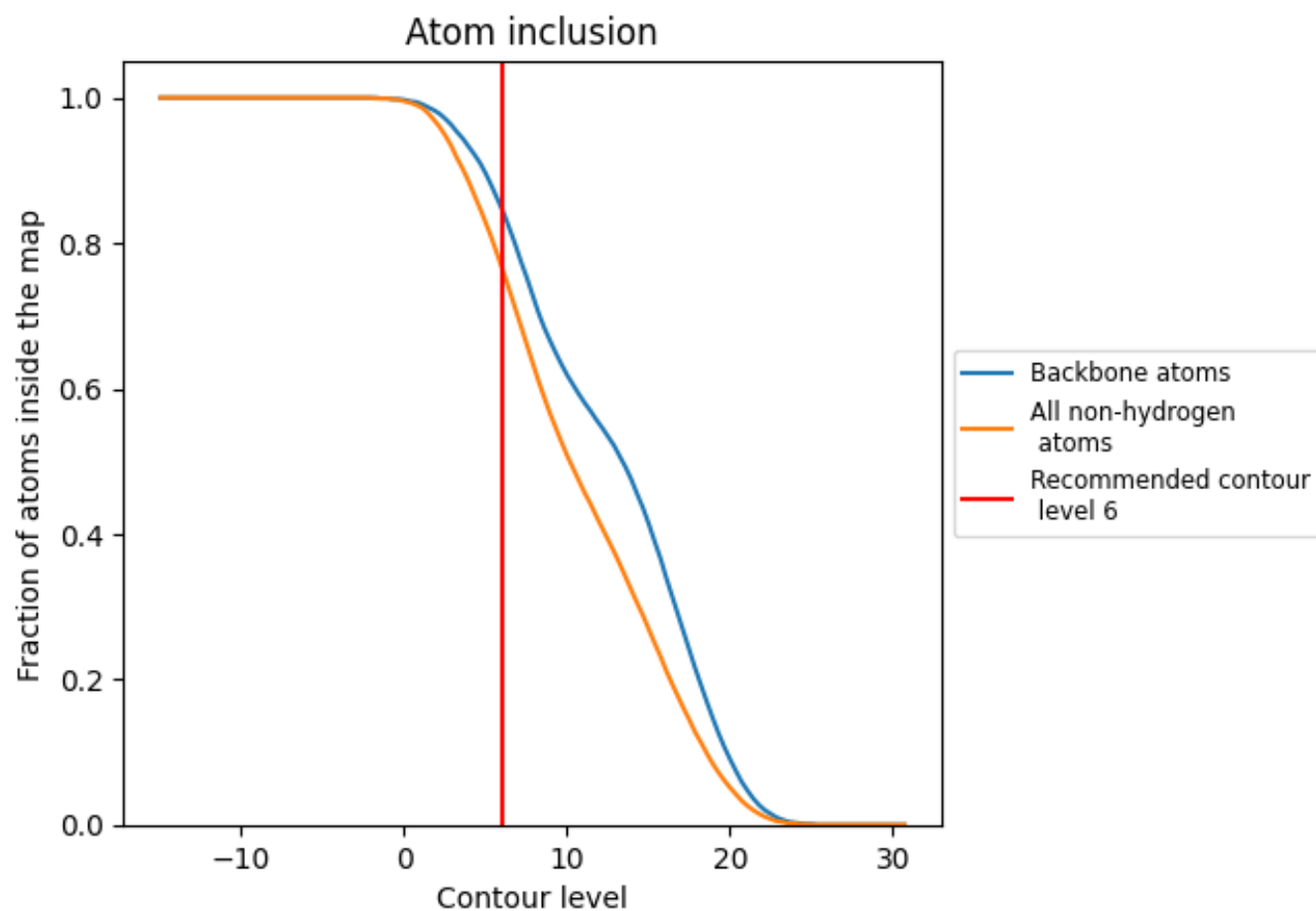


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 85% 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 (6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7680	 0.5520
A	 0.8410	 0.5580
B	 0.7920	 0.5500
C	 0.7900	 0.5300
D	 0.7660	 0.5440
E	 0.8440	 0.5670
F	 0.8200	 0.5590
G	 0.8240	 0.5580
H	 0.7700	 0.5600
I	 0.8840	 0.5870
J	 0.8850	 0.5910
K	 0.8920	 0.5860
L	 0.8650	 0.5830
M	 0.8790	 0.5840
N	 0.8190	 0.5710
O	 0.8270	 0.5520
P	 0.7840	 0.5440
Q	 0.7940	 0.5290
R	 0.7710	 0.5420
S	 0.8380	 0.5610
T	 0.8230	 0.5570
U	 0.8180	 0.5530
V	 0.7680	 0.5520
W	 0.8760	 0.5860
X	 0.8750	 0.5850
Y	 0.8920	 0.5870
Z	 0.8720	 0.5860
a	 0.8870	 0.5820
b	 0.8120	 0.5690
c	 0.3640	 0.4750
d	 0.5550	 0.5160
e	 0.4520	 0.5370
f	 0.3660	 0.4500
g	 0.5560	 0.5080
h	 0.4490	 0.5330

