



wwPDB EM Validation Summary Report ⓘ

Aug 10, 2026 – 03:11 PM EDT

PDB ID : 35ZW / pdb_000035zw
EMDB ID : EMD-77308
Title : yeast 26S proteasome base assembly intermediate, Hsm3-Rpt1-Rpt2 (base-Hsm3-Nas6)
Authors : Hsieh, H.H.; Martin, A.
Deposited on : 2026-05-26
Resolution : 3.31 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

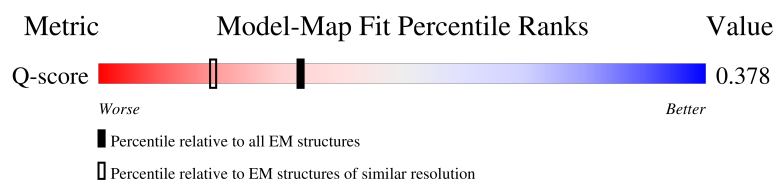
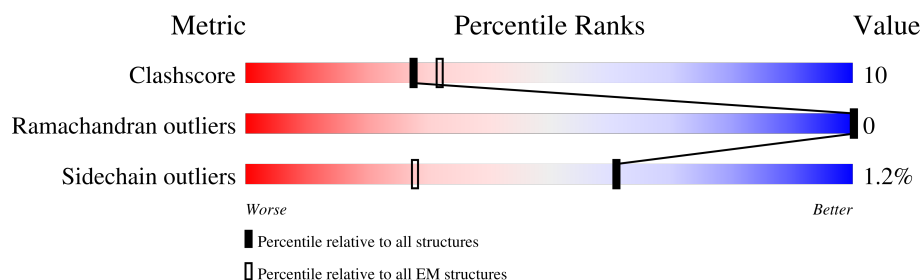
EMDB validation analysis : 0.0.1.dev133
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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) : 2.50

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 3.31 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.



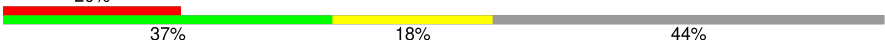
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14550 (2.81 - 3.81)

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	467	
2	B	437	
3	D	480	
4	G	437	

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Mol	Chain	Length	Quality of chain
5	H	434	 50% 10% 39%
6	L	428	 20% 37% 18% 44%
7	O	228	 98% 75% 25%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 15764 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 26S proteasome regulatory subunit 7 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	253	Total	C	N	O	S	0	0
			1961	1231	352	365	13		

- Molecule 2 is a protein called 26S proteasome regulatory subunit 4 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	256	Total	C	N	O	S	0	0
			1991	1252	342	385	12		

- Molecule 3 is a protein called DNA mismatch repair protein HSM3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	465	Total	C	N	O	S	0	0
			3805	2454	608	730	13		

- Molecule 4 is a protein called 26S proteasome subunit RPT4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	273	Total	C	N	O	S	0	0
			2133	1353	369	401	10		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	282	GLN	GLU	engineered mutation	UNP P53549

- Molecule 5 is a protein called 26S proteasome regulatory subunit 6A.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	266	Total	C	N	O	S	0	0
			2062	1287	364	402	9		

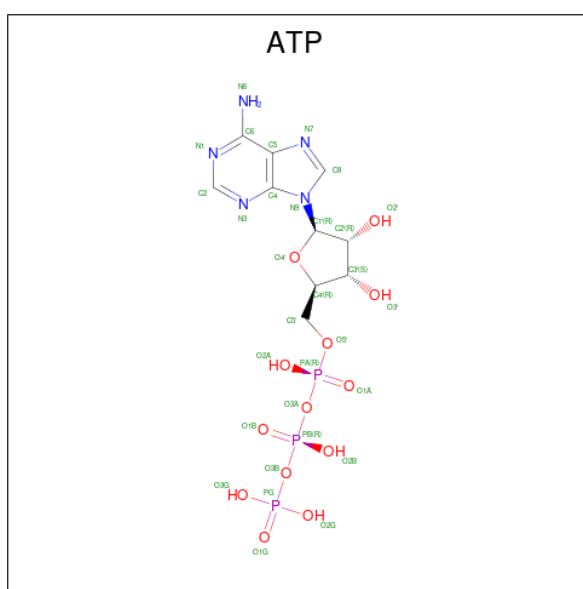
- Molecule 6 is a protein called 26S proteasome regulatory subunit 6B homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L	238	Total	C	N	O	S	0	0
			1849	1153	334	354	8		

- Molecule 7 is a protein called Probable 26S proteasome regulatory subunit p28.

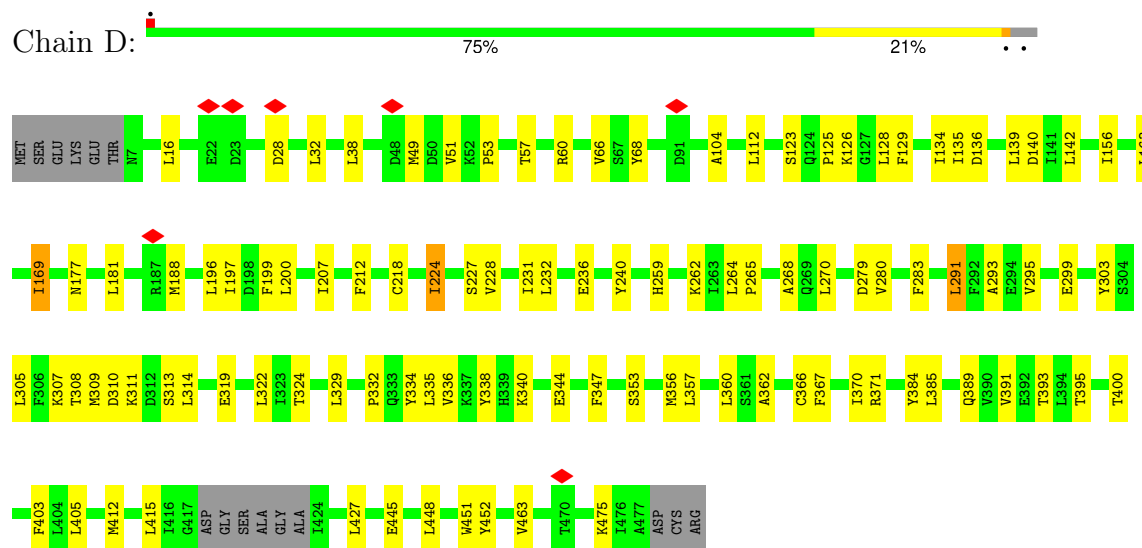
Mol	Chain	Residues	Atoms					AltConf	Trace
7	O	228	Total	C	N	O	S	0	0
			1808	1159	307	335	7		

- Molecule 8 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).

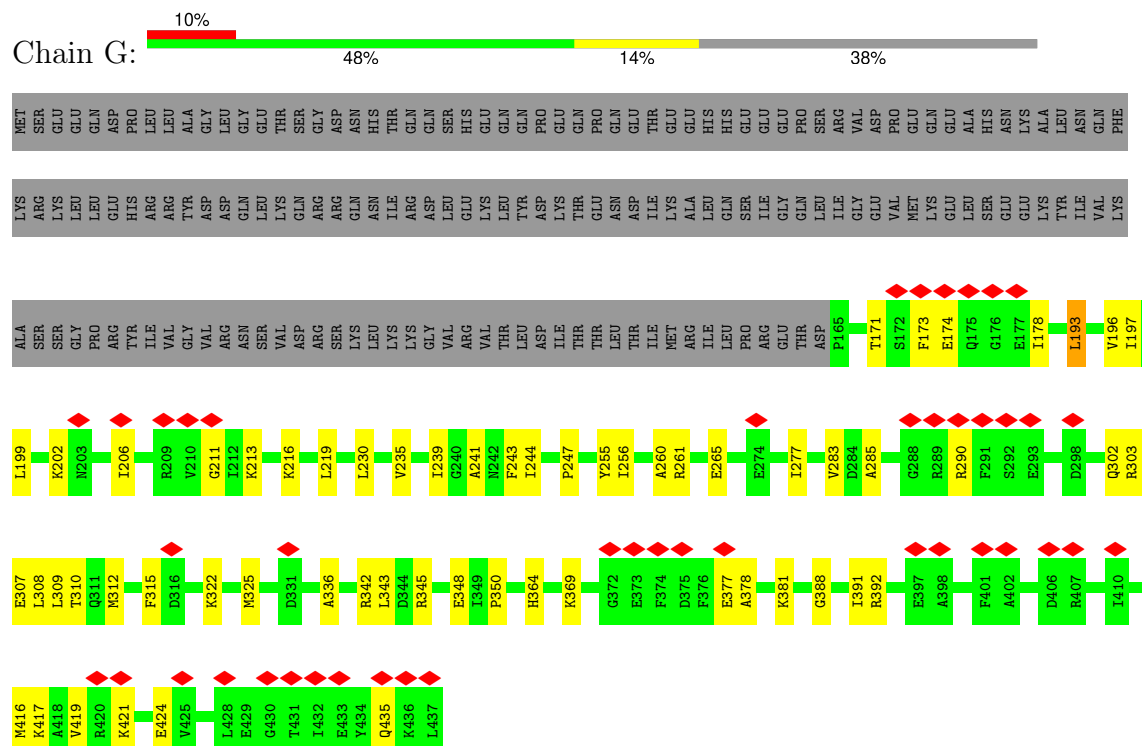


Mol	Chain	Residues	Atoms					AltConf
8	A	1	Total	C	N	O	P	0
			31	10	5	13	3	
8	B	1	Total	C	N	O	P	0
			31	10	5	13	3	
8	G	1	Total	C	N	O	P	0
			31	10	5	13	3	
8	H	1	Total	C	N	O	P	0
			31	10	5	13	3	
8	L	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 3: DNA mismatch repair protein HSM3

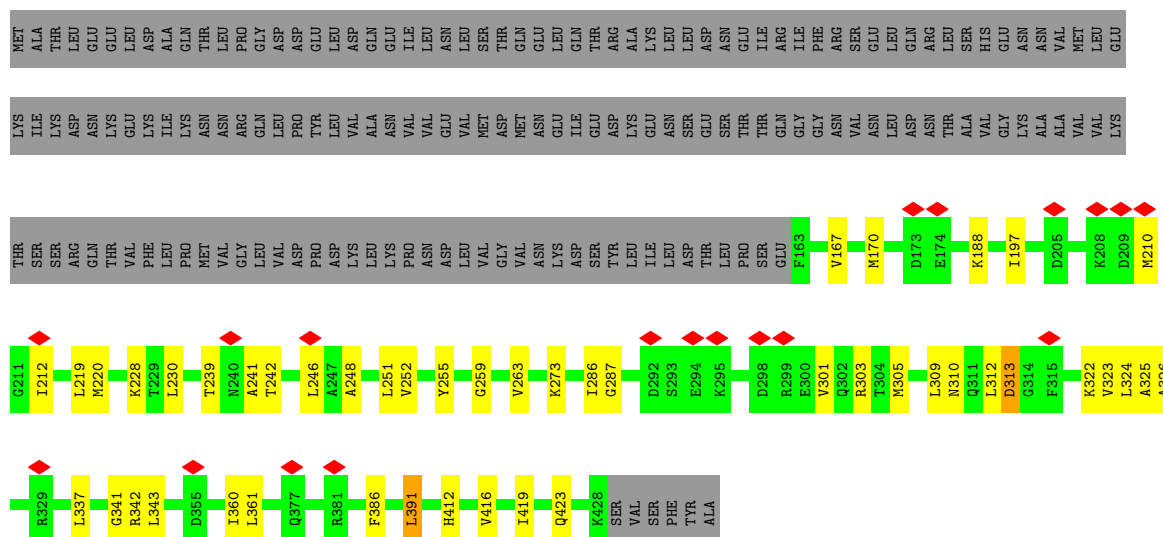


- Molecule 4: 26S proteasome subunit RPT4

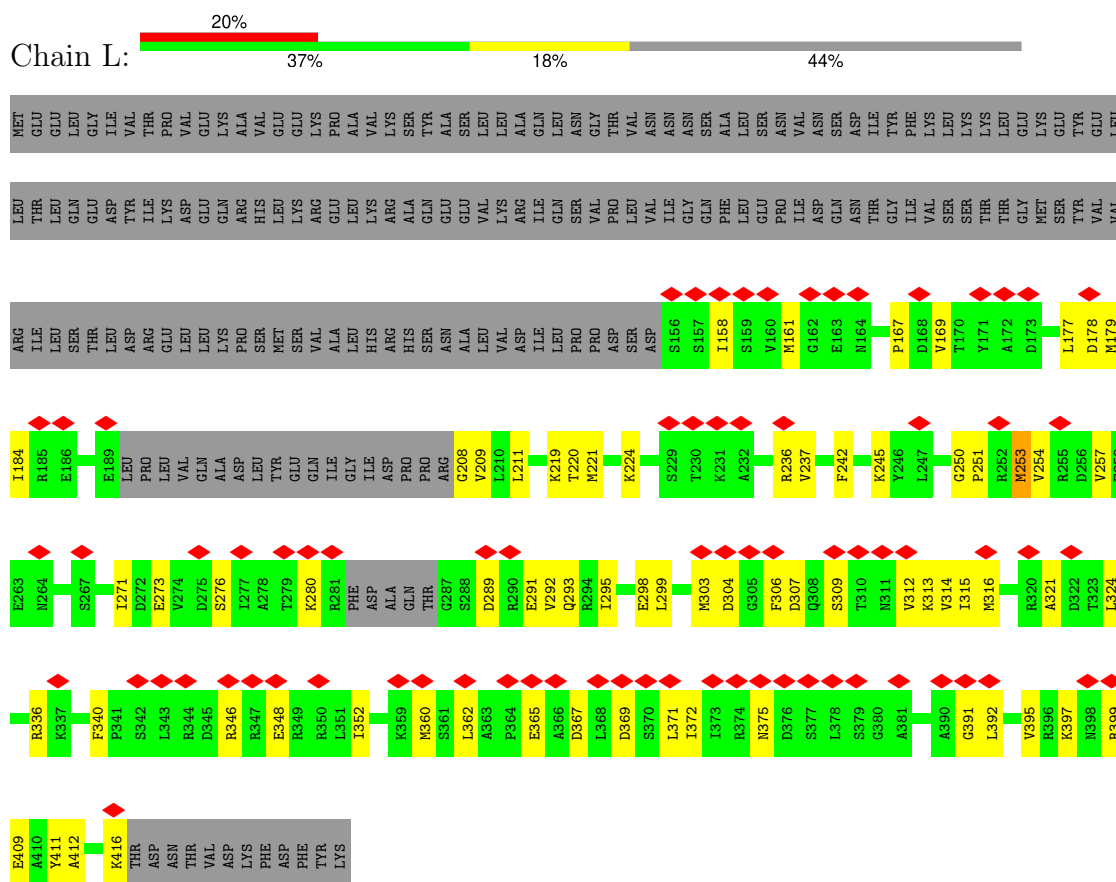


- Molecule 5: 26S proteasome regulatory subunit 6A

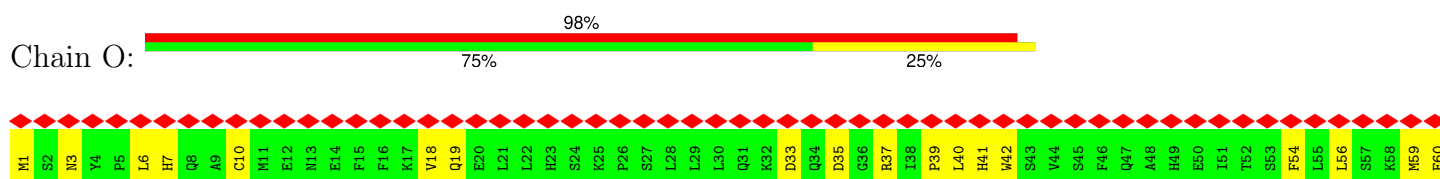




• Molecule 6: 26S proteasome regulatory subunit 6B homolog



• Molecule 7: Probable 26S proteasome regulatory subunit p28



A181	L182	A183	E184	G185	H186	G187	D188	A189	A190	V191	L192	L193	V194	E195	K196	Y197	G198	A199	E200	Y201	D202	L203	V204	D205	N206	K207	G208	A209	K210	A211	E212	D213	V214	A215	L216	N217	E218	Q219	V220	K221	K222	F223	F224	L225	N226	N227	V228												
E121	V122	S123	Q124	F125	L126	I127	E128	N129	G130	A131	S132	V133	R134	I135	K136	D137	K138	F139	N140	Q141	I142	P143	L144	H145	R146	A147	A148	S149	V150	G151	S152	L153	K154	L155	I156	E157	L158	L159	C160	G161	L162	G163	K164	S165	A166	V167	N168	W169	Q170	D171	K172	Q173	G174	W175	T176	P177	L178	F179	H180
N61	V62	N63	L64	D65	D66	Y67	P68	D69	D70	S71	G72	W73	T74	P75	F76	H77	I78	A79	C80	S81	V82	G83	N84	L85	E86	V87	V88	K89	S90	L91	Y92	D93	R94	P95	L96	K97	P98	D99	L100	N101	K102	I103	T104	N105	Q106	G107	V108	T109	C110	L111	H112	L113	A114	V115	G116	K117	K118	W119	F120

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	241263	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.646	Depositor
Minimum map value	-0.310	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.11	Depositor
Map size (Å)	536.576, 536.576, 536.576	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.048, 1.048, 1.048	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ATP

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	0.15	0/1990	0.32	0/2679
2	B	0.15	0/2014	0.41	0/2708
3	D	0.16	0/3872	0.37	0/5243
4	G	0.17	0/2170	0.40	0/2921
5	H	0.16	0/2088	0.38	0/2809
6	L	0.15	0/1868	0.40	0/2513
7	O	0.14	0/1851	0.40	0/2506
All	All	0.16	0/15853	0.38	0/21379

There are no bond length outliers.

There are no bond angle outliers.

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	1961	0	2004	27	0
2	B	1991	0	2067	43	0
3	D	3805	0	3842	73	0
4	G	2133	0	2165	50	0
5	H	2062	0	2095	34	0
6	L	1849	0	1897	60	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	O	1808	0	1792	40	0
8	A	31	0	12	0	0
8	B	31	0	12	0	0
8	G	31	0	12	3	0
8	H	31	0	12	2	0
8	L	31	0	12	2	0
All	All	15764	0	15922	306	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 10.

The worst 5 of 306 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:H:361:LEU:HD13	5:H:391:LEU:HG	1.65	0.78
7:O:64:LEU:HD11	7:O:91:LEU:HD22	1.69	0.74
4:G:219:LEU:HB2	4:G:343:LEU:HD13	1.71	0.73
6:L:330:ARG:HH11	6:L:333:ARG:HG3	1.53	0.73
3:D:336:VAL:HG21	3:D:366:CYS:HA	1.71	0.72

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	249/467 (53%)	246 (99%)	3 (1%)	0	100	100
2	B	252/437 (58%)	245 (97%)	7 (3%)	0	100	100
3	D	461/480 (96%)	451 (98%)	10 (2%)	0	100	100
4	G	271/437 (62%)	267 (98%)	4 (2%)	0	100	100
5	H	264/434 (61%)	256 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	L	232/428 (54%)	226 (97%)	6 (3%)	0	100	100
7	O	226/228 (99%)	219 (97%)	7 (3%)	0	100	100
All	All	1955/2911 (67%)	1910 (98%)	45 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	211/399 (53%)	210 (100%)	1 (0%)	81	82
2	B	222/385 (58%)	217 (98%)	5 (2%)	44	66
3	D	434/445 (98%)	428 (99%)	6 (1%)	59	72
4	G	226/377 (60%)	225 (100%)	1 (0%)	84	83
5	H	221/375 (59%)	218 (99%)	3 (1%)	59	72
6	L	200/374 (54%)	197 (98%)	3 (2%)	57	71
7	O	198/198 (100%)	197 (100%)	1 (0%)	81	82
All	All	1712/2553 (67%)	1692 (99%)	20 (1%)	61	74

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	H	313	ASP
6	L	306	PHE
7	O	224	PHE
6	L	313	LYS
3	D	28	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
4	G	311	GLN
7	O	105	ASN
5	H	328	ASN
7	O	173	GLN
6	L	404	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](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
8	ATP	L	501	-	32,33,33	0.32	0	48,52,52	0.35	0
8	ATP	G	501	-	32,33,33	0.45	0	48,52,52	0.38	0
8	ATP	A	501	-	32,33,33	0.44	0	48,52,52	0.34	0
8	ATP	H	501	-	32,33,33	0.37	0	48,52,52	0.34	0
8	ATP	B	501	-	32,33,33	0.33	0	48,52,52	0.34	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	ATP	L	501	-	-	2/22/38/38	0/3/3/3
8	ATP	G	501	-	-	5/22/38/38	0/3/3/3
8	ATP	A	501	-	-	6/22/38/38	0/3/3/3
8	ATP	H	501	-	-	5/22/38/38	0/3/3/3
8	ATP	B	501	-	-	4/22/38/38	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

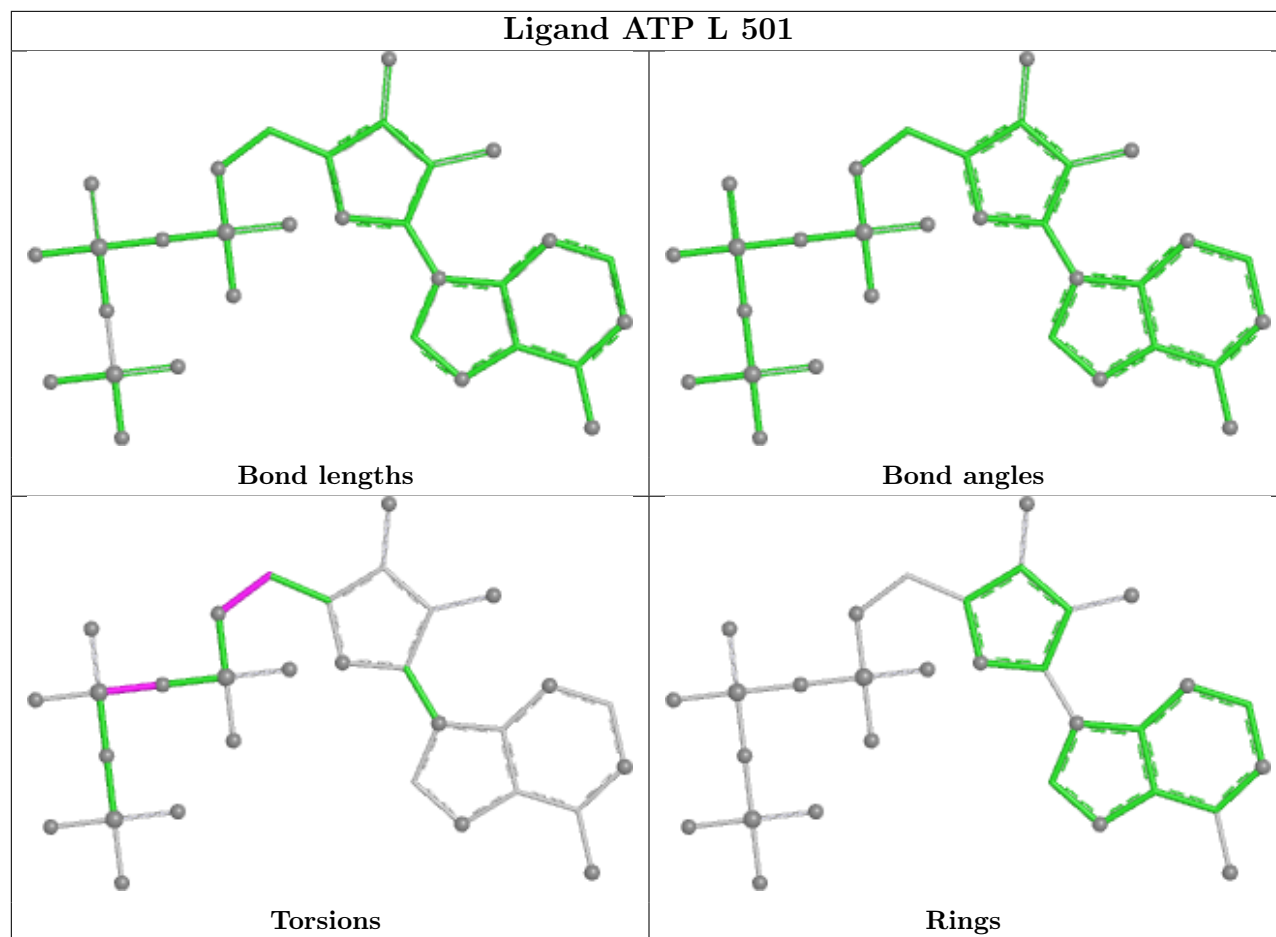
Mol	Chain	Res	Type	Atoms
8	B	501	ATP	PB-O3B-PG-O2G
8	B	501	ATP	C5'-O5'-PA-O2A
8	G	501	ATP	C5'-O5'-PA-O1A
8	G	501	ATP	C5'-O5'-PA-O3A
8	G	501	ATP	C4'-C5'-O5'-PA

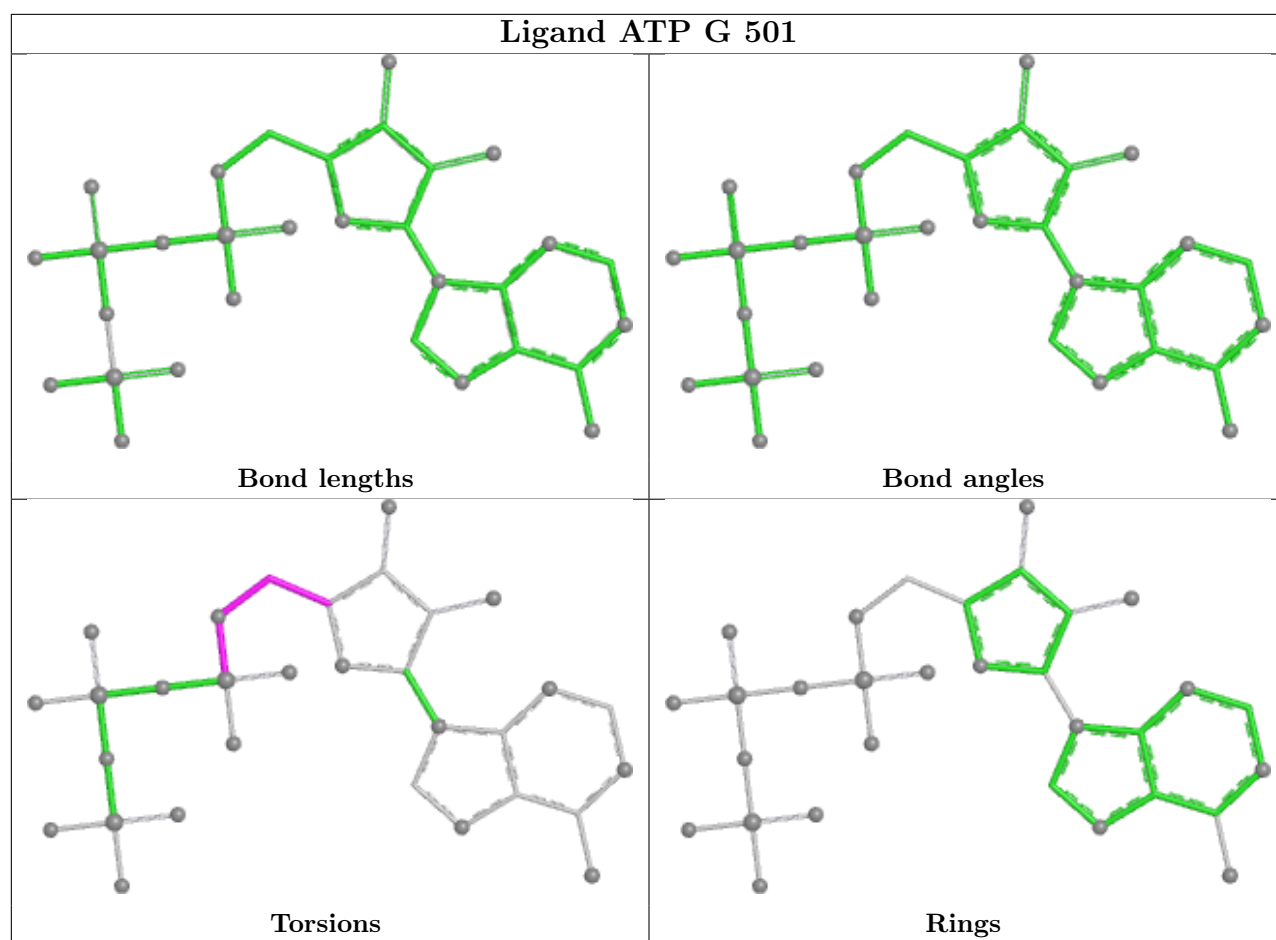
There are no ring outliers.

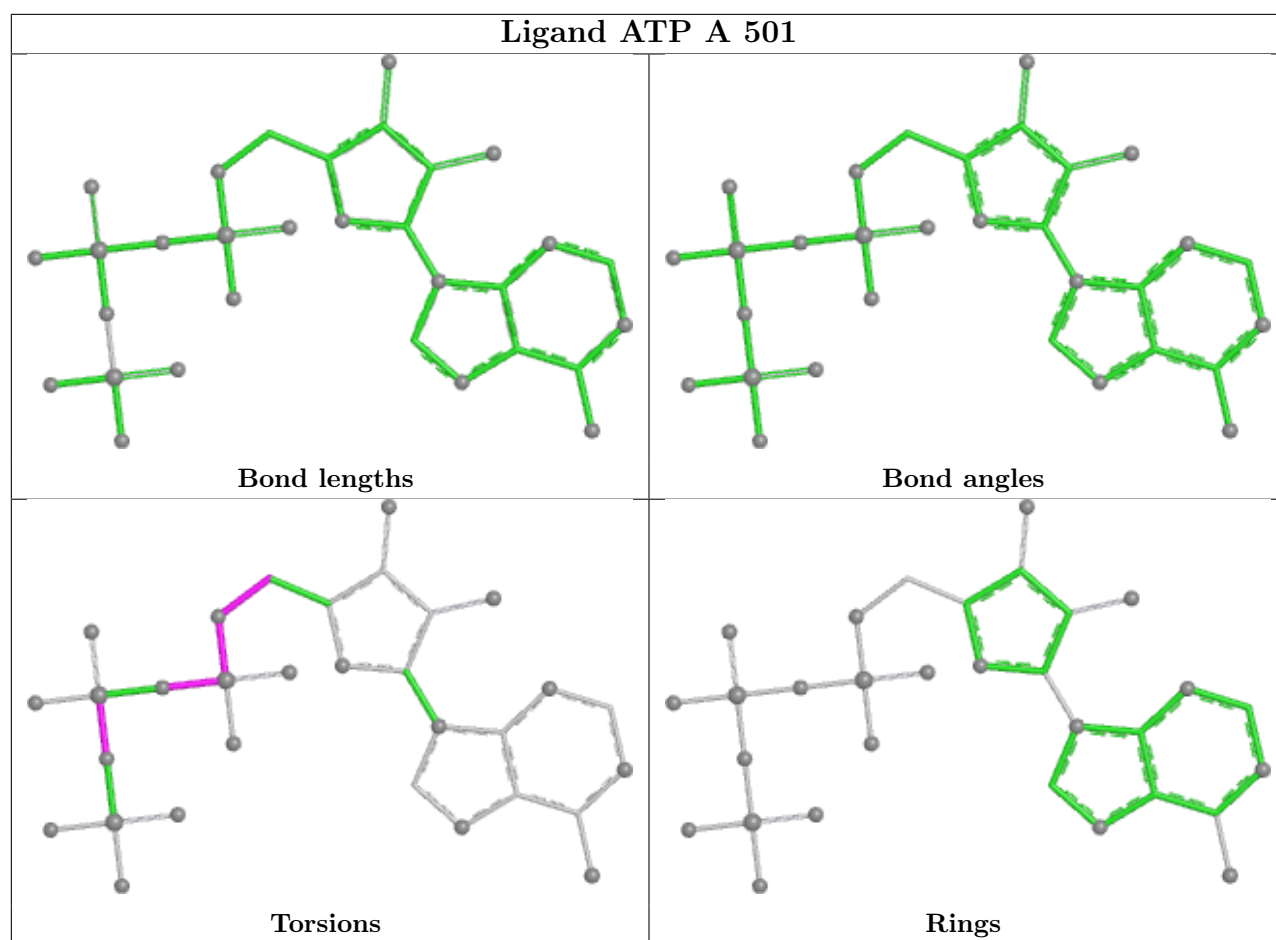
3 monomers are involved in 7 short contacts:

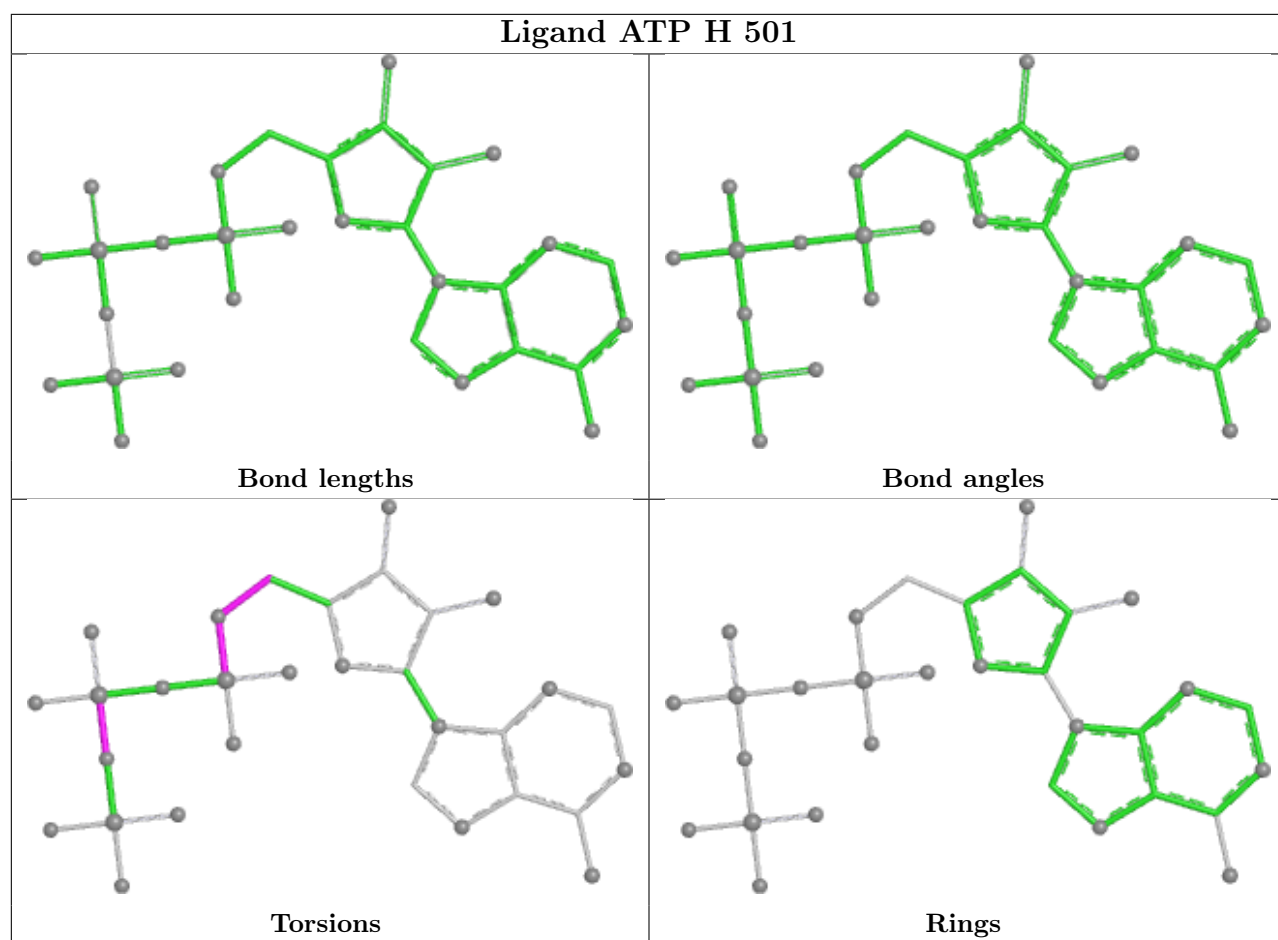
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	L	501	ATP	2	0
8	G	501	ATP	3	0
8	H	501	ATP	2	0

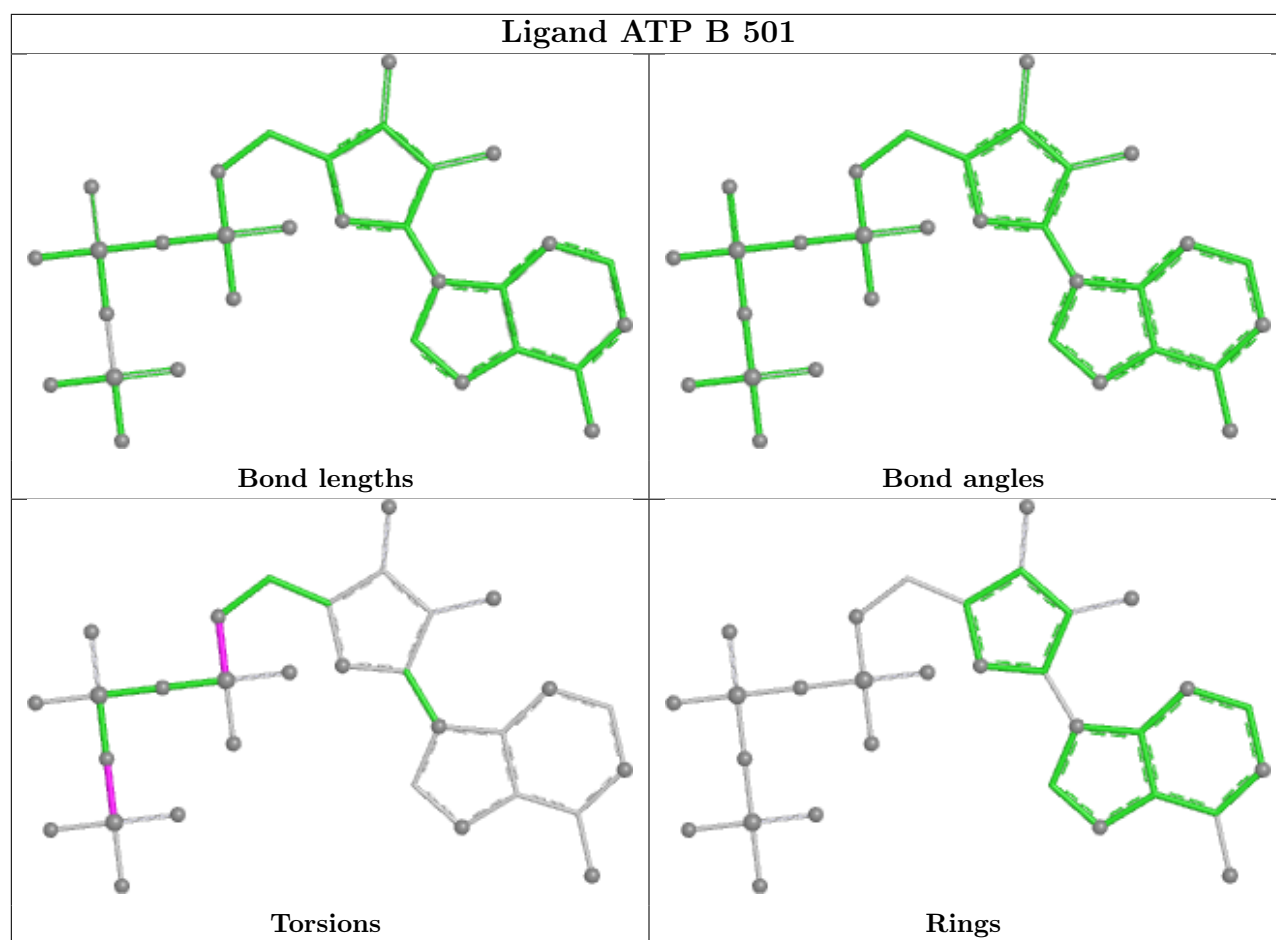
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











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.

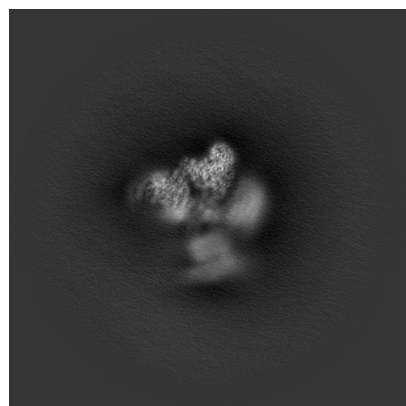
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-77308. These allow visual inspection of the internal detail of the map and identification of artifacts.

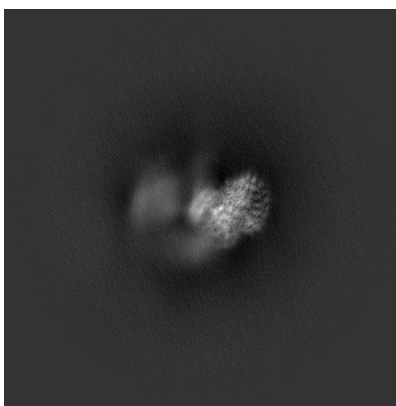
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

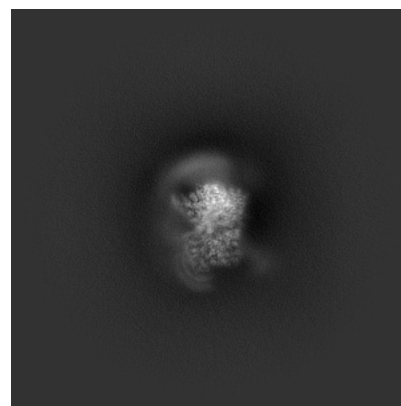
6.1.1 Primary map



X

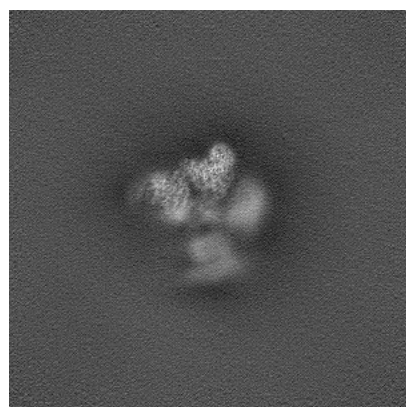


Y

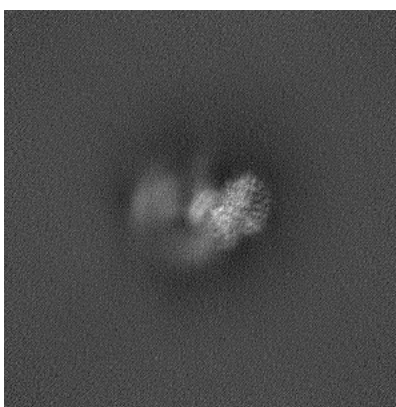


Z

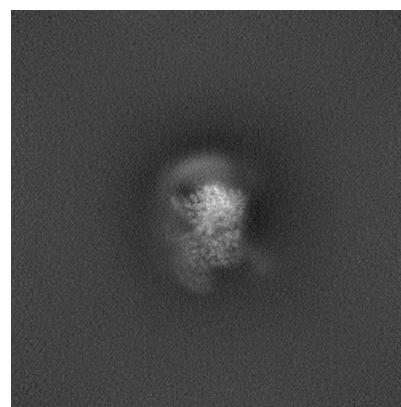
6.1.2 Raw map



X



Y



Z

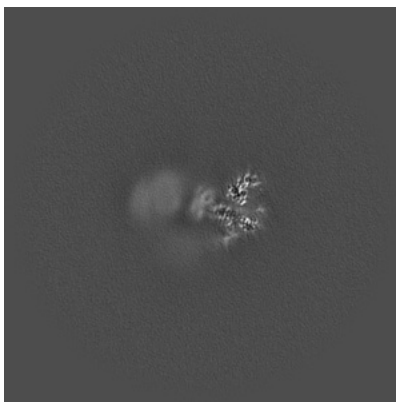
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

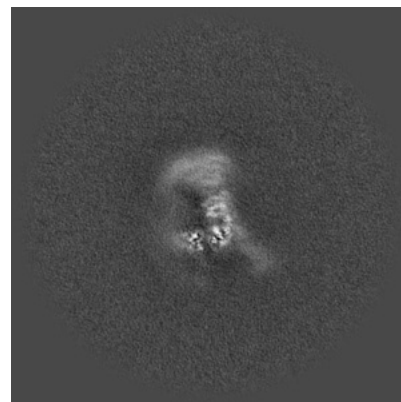
6.2.1 Primary map



X Index: 256



Y Index: 256



Z Index: 256

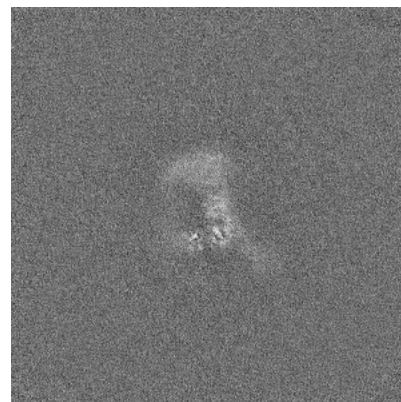
6.2.2 Raw map



X Index: 256



Y Index: 256



Z Index: 256

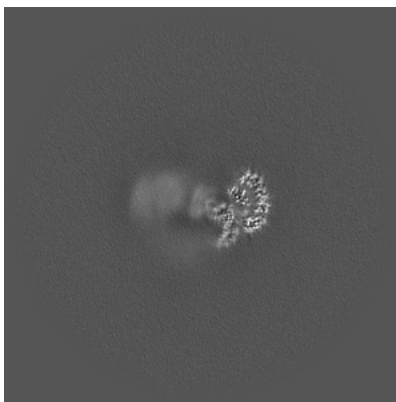
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

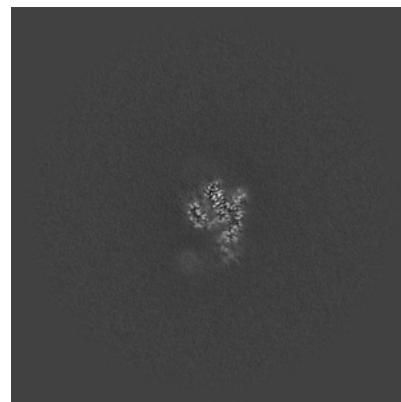
6.3.1 Primary map



X Index: 260

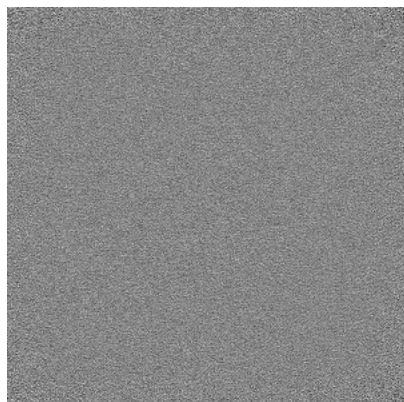


Y Index: 263

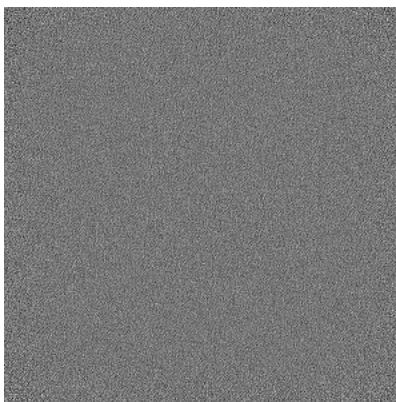


Z Index: 304

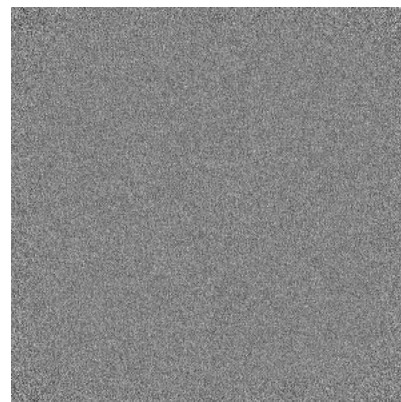
6.3.2 Raw map



X Index: 0



Y Index: 0

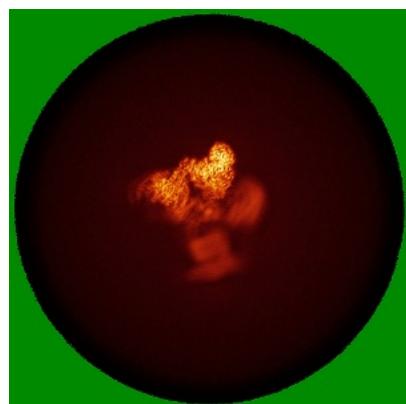


Z Index: 511

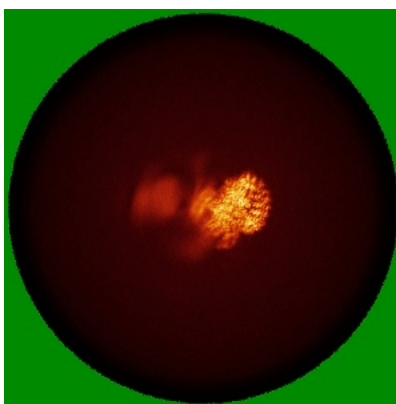
The images above show the largest variance slices of the map in three orthogonal directions.

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

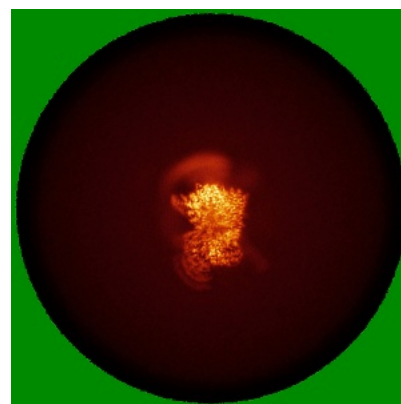
6.4.1 Primary map



X

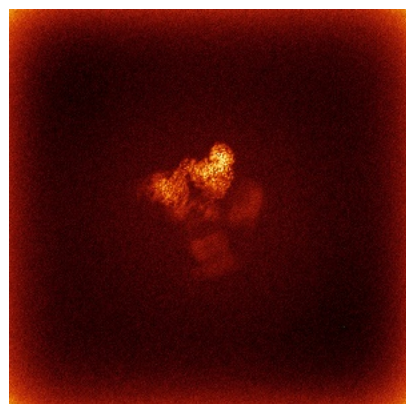


Y

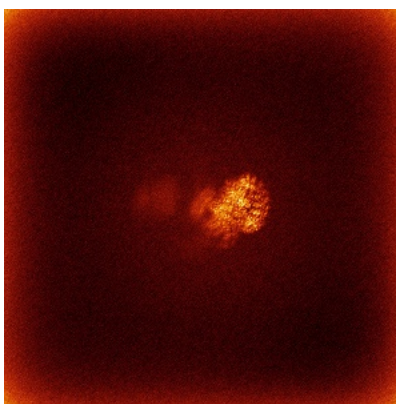


Z

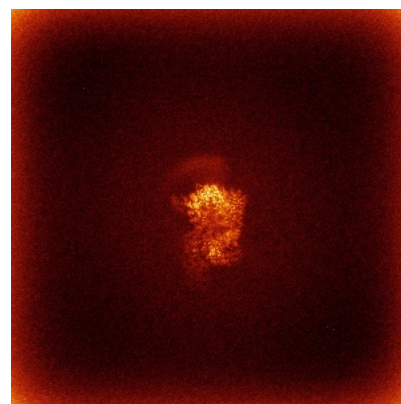
6.4.2 Raw map



X



Y

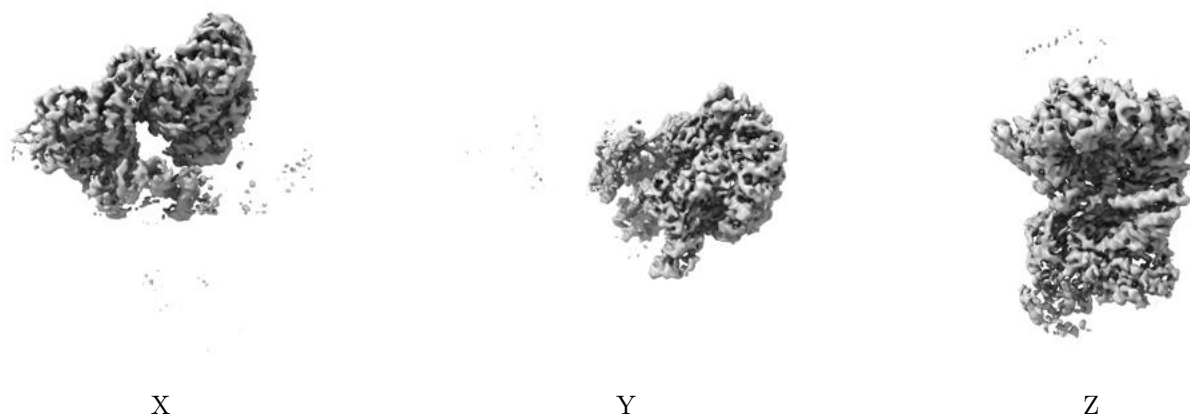


Z

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.

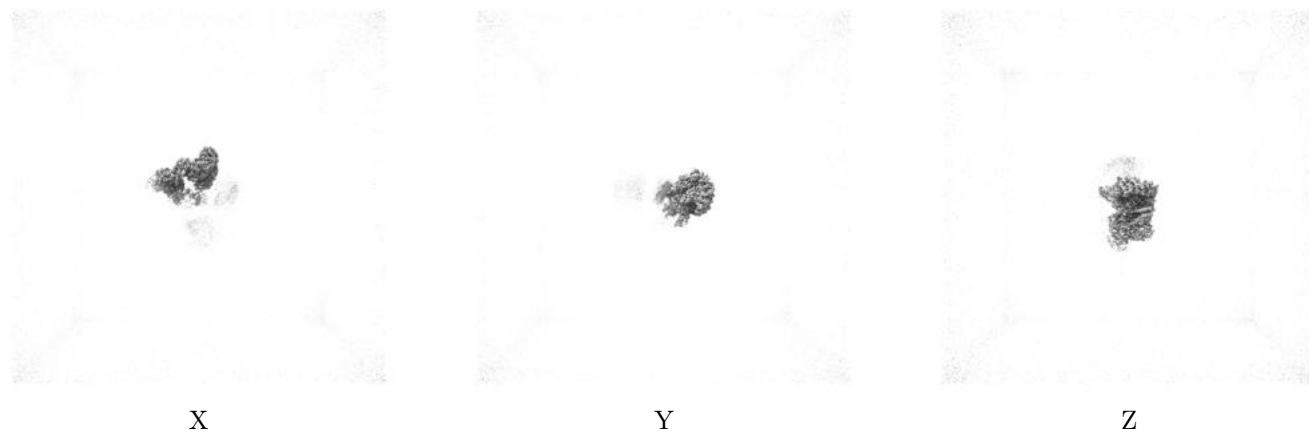
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.11. 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.

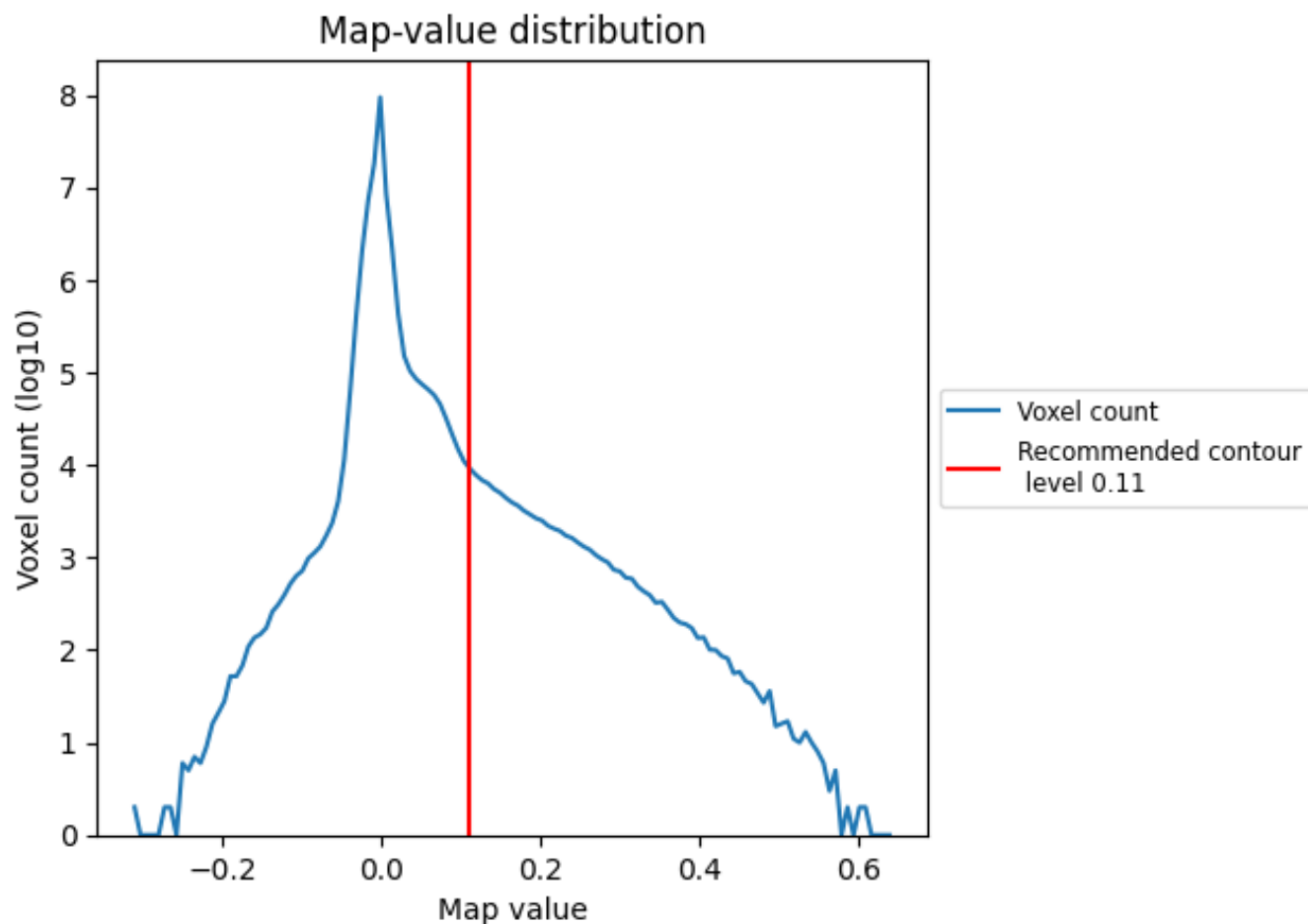
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

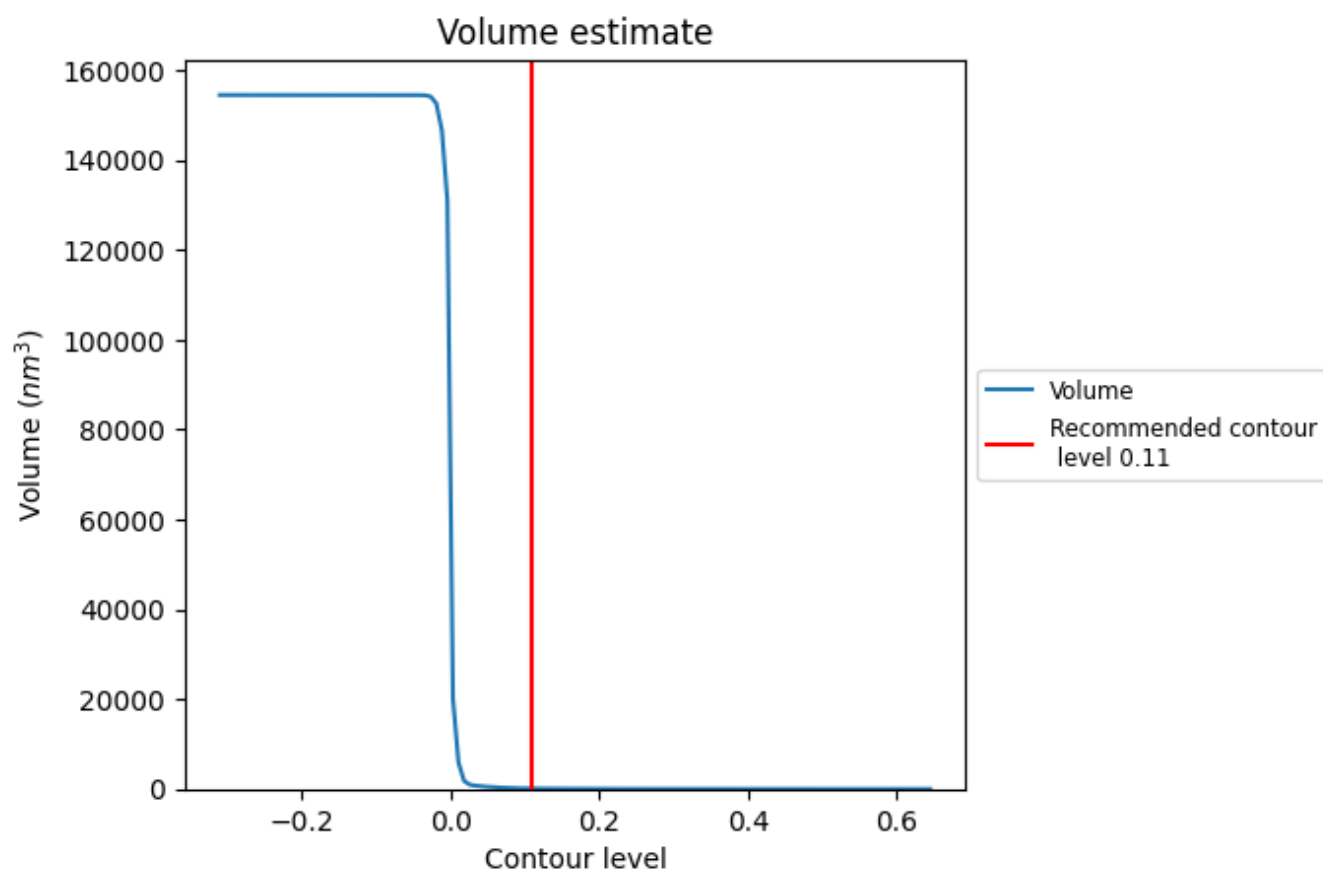
This section contains the results of statistical analysis of the map.

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.

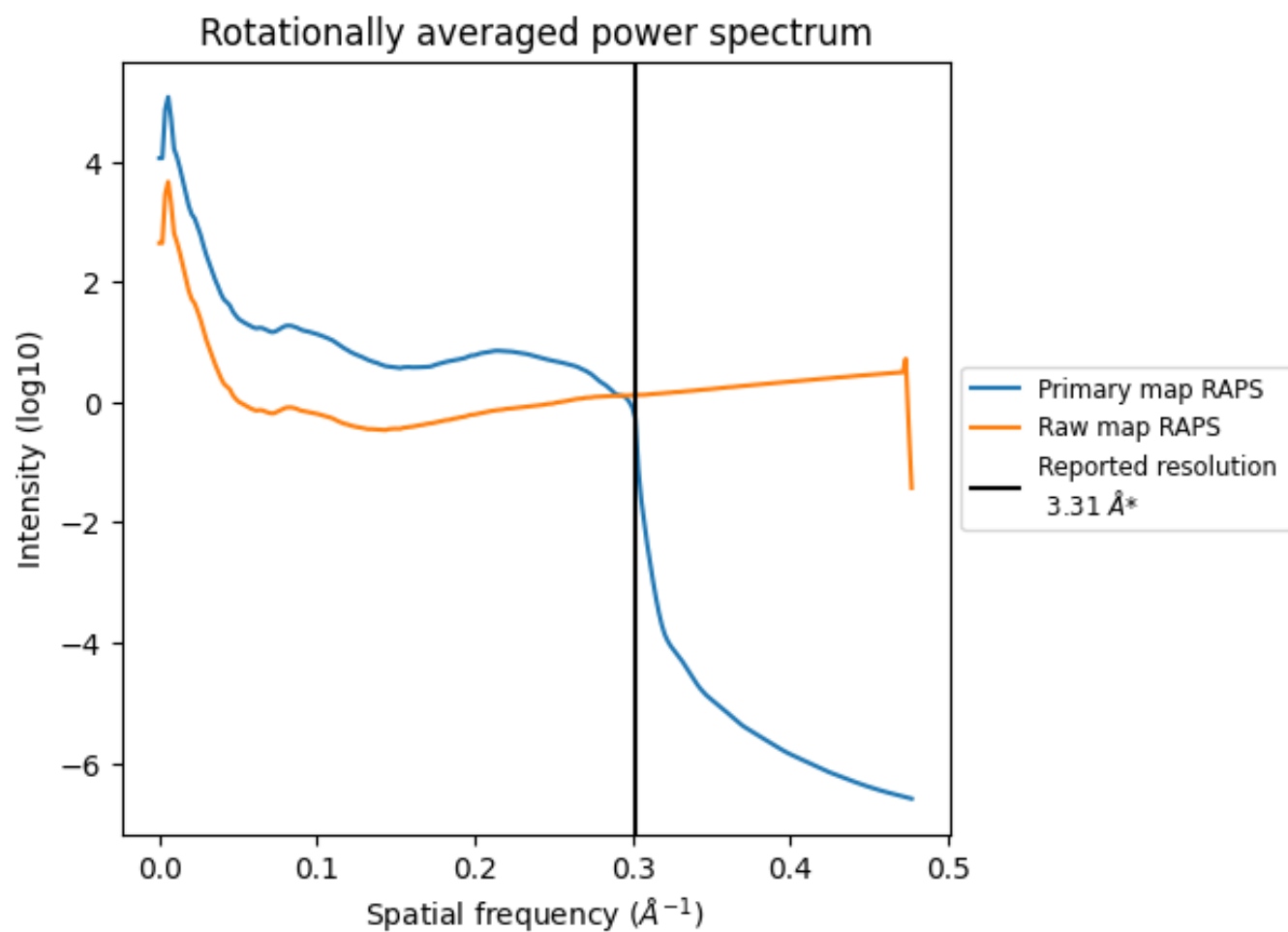
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 102 nm³; this corresponds to an approximate mass of 92 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 ⓘ

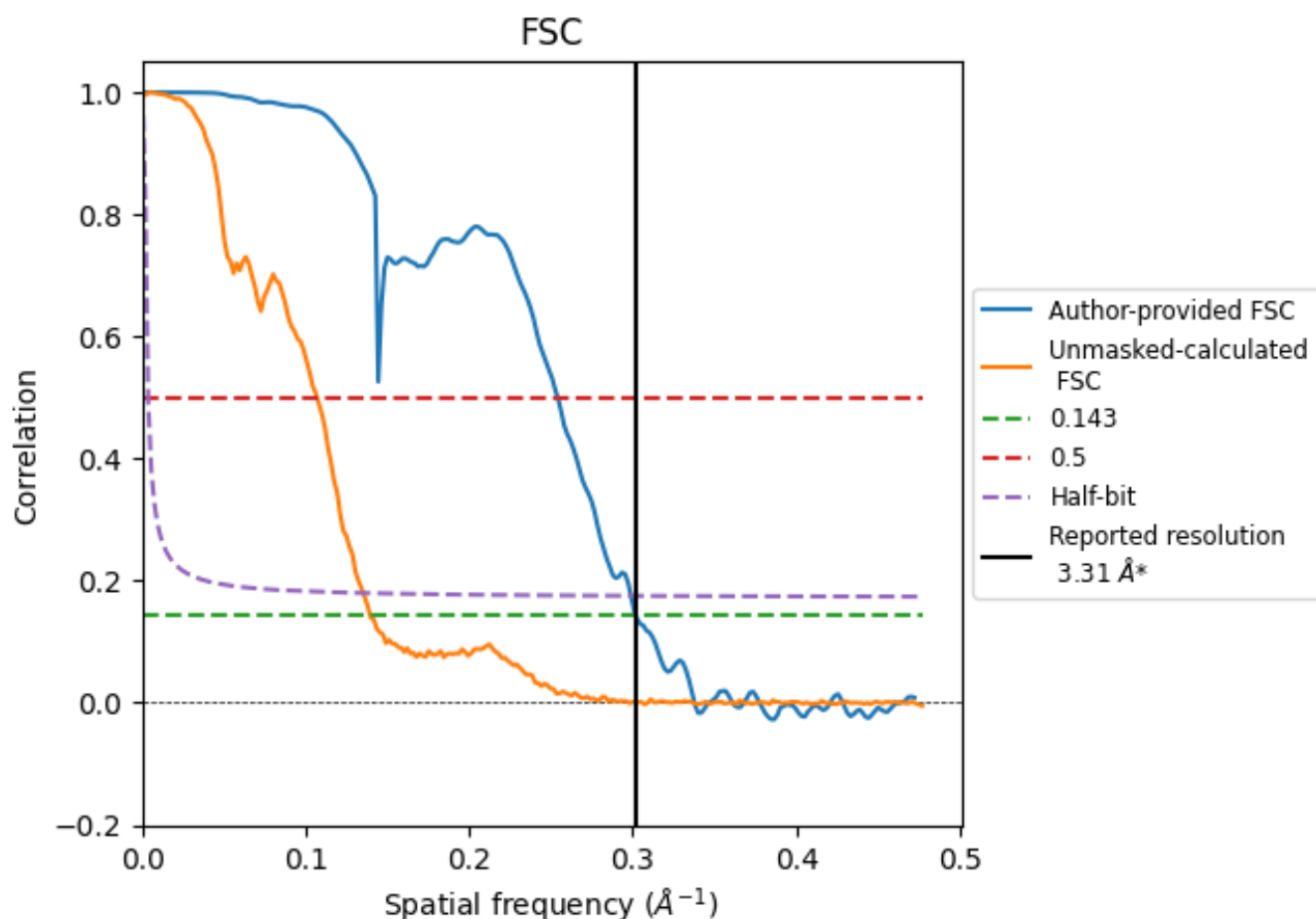


*Reported resolution corresponds to spatial frequency of $0.302 \text{ \AA}^{-1}$

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 0.302 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

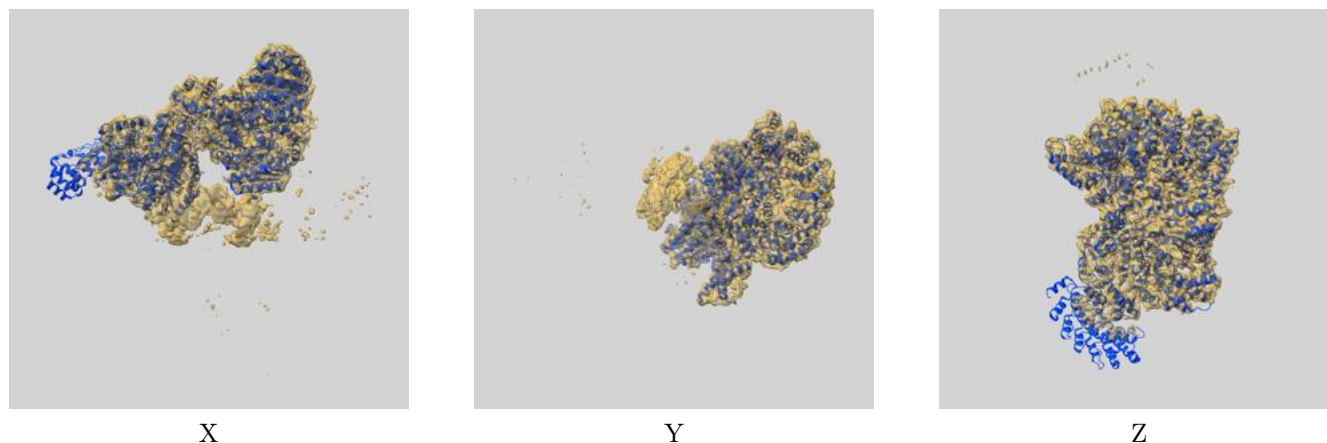
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.31	-	-
Author-provided FSC curve	3.31	3.93	3.34
Unmasked-calculated*	7.17	9.35	7.40

*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 7.17 differs from the reported value 3.31 by more than 10 %

9 Map-model fit [i](#)

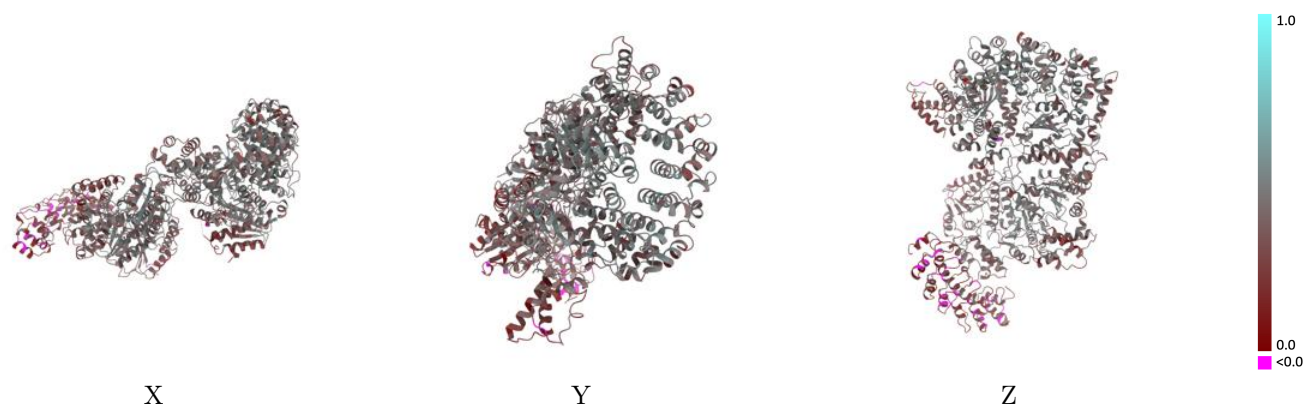
This section contains information regarding the fit between EMDB map EMD-77308 and PDB model 35ZW. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)



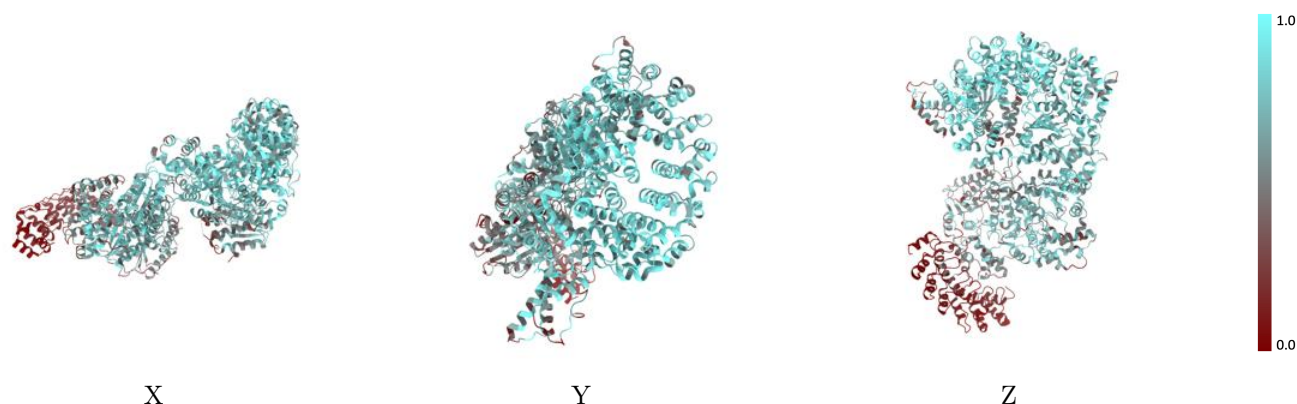
The images above show the 3D surface view of the map at the recommended contour level 0.11 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



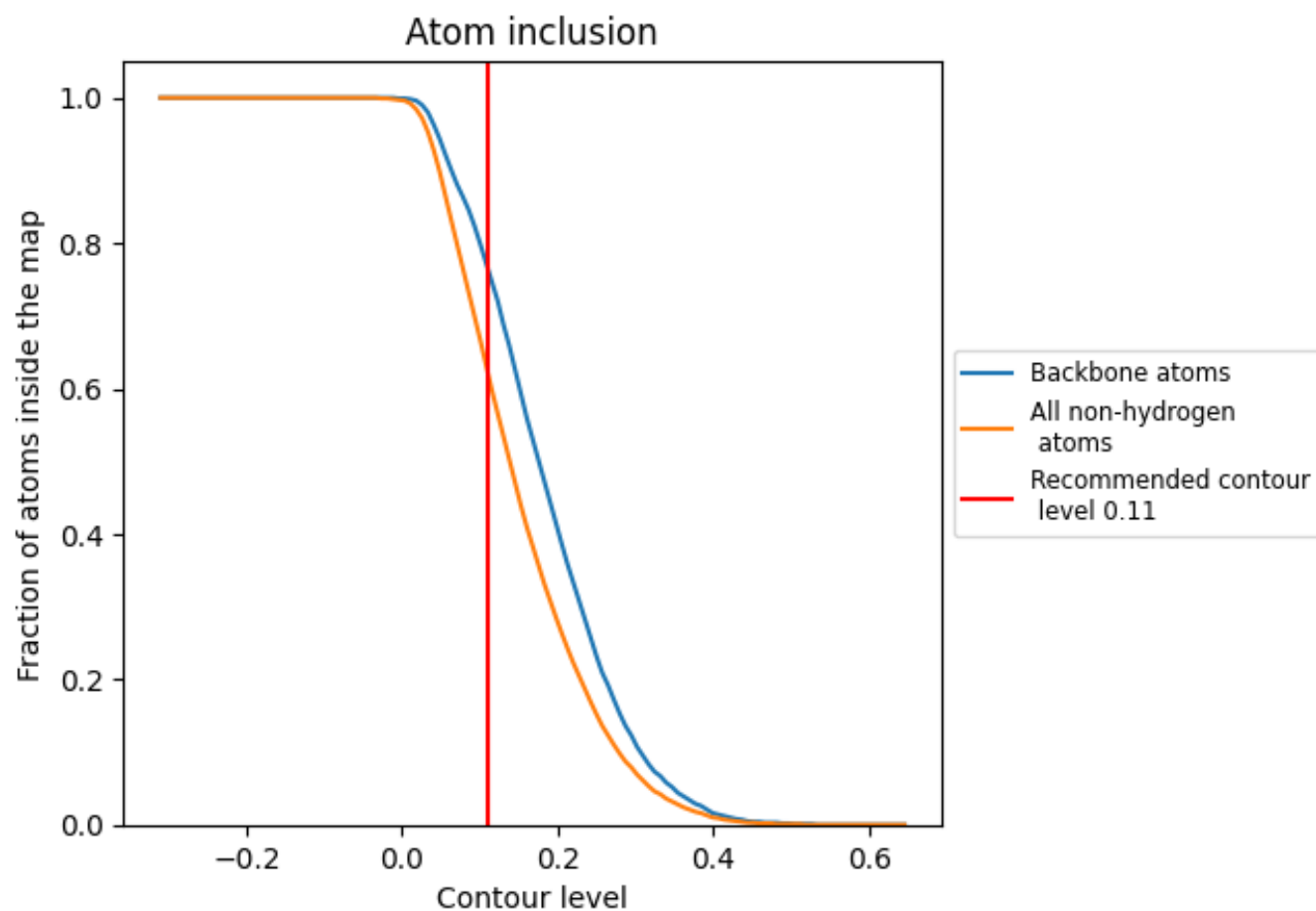
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.11).

9.4 Atom inclusion [i](#)



At the recommended contour level, 77% 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 (0.11) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6230	0.3780
A	0.8300	0.4580
B	0.6150	0.3450
D	0.8050	0.4410
G	0.6630	0.4000
H	0.6920	0.4160
L	0.4890	0.3250
O	0.0350	0.1760

1.0

0.0

<0.0