



wwPDB EM Validation Summary Report ⓘ

Aug 6, 2026 – 12:48 AM JST

PDB ID : 5Z3U / pdb_00005z3u
EMDB ID : EMD-6882
Title : Structure of Snf2-nucleosome complex at shl2 in ADP BeFx state
Authors : Li, M.; Xia, X.; Liu, X.; Li, X.; Chen, Z.
Deposited on : 2018-01-08
Resolution : 4.31 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:

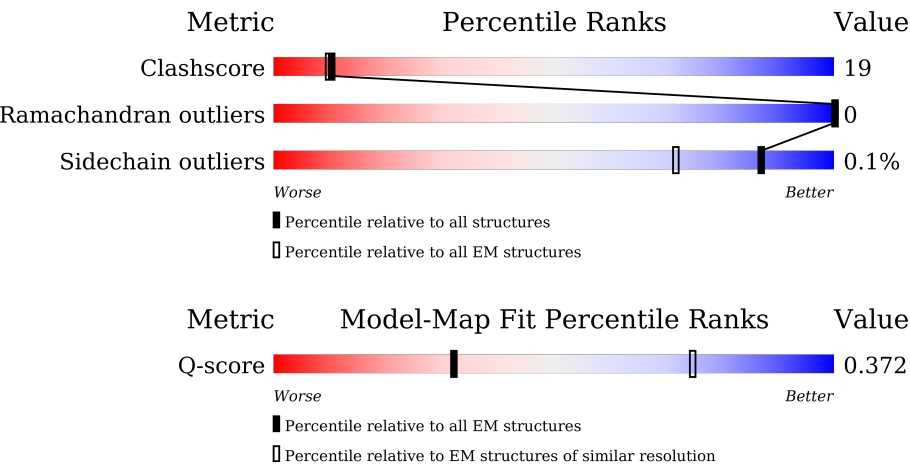
EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using 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 i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.31 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.



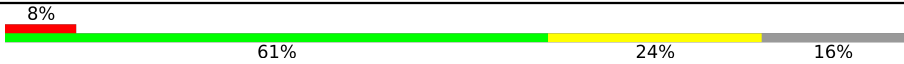
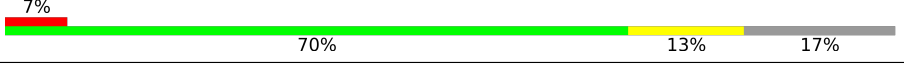
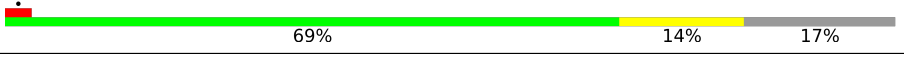
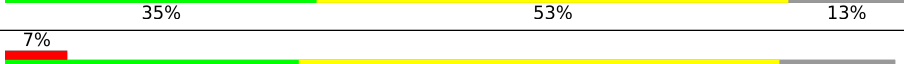
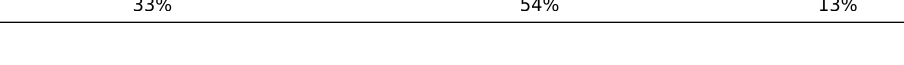
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	3904 (3.81 - 4.80)

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	O	735	50% 34% 44% 22%
2	A	135	50% 22% 27%

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Mol	Chain	Length	Quality of chain
3	F	102	
4	C	129	
4	G	129	
5	D	122	
5	H	122	
6	I	167	
7	J	167	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	BEF	O	1503	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 16766 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 Transcription regulatory protein SNF2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	O	575	Total	C	N	O	S	0	0
			4723	3004	831	871	17		

- Molecule 2 is a protein called Histone H3.2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	98	Total	C	N	O	S	0	0
			801	506	153	139	3		
2	E	95	Total	C	N	O	S	0	0
			779	492	148	136	3		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	GLY	conflict	UNP P84233
E	102	ALA	GLY	conflict	UNP P84233

- Molecule 3 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	87	Total	C	N	O	S	0	0
			703	443	142	117	1		
3	F	86	Total	C	N	O	S	0	0
			672	424	130	117	1		

- Molecule 4 is a protein called Histone H2A.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	C	107	Total	C	N	O	0	0
			811	510	158	143		
4	G	107	Total	C	N	O	0	0
			815	513	159	143		

- Molecule 5 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	D	93	Total	C	N	O	S	0	0
			718	451	128	137	2		
5	H	93	Total	C	N	O	S	0	0
			726	457	130	137	2		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	29	THR	SER	see sequence details	UNP P02281
H	29	THR	SER	see sequence details	UNP P02281

- Molecule 6 is a DNA chain called DNA (167-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	I	146	Total	C	N	O	P	0	0
			2975	1413	540	876	146		

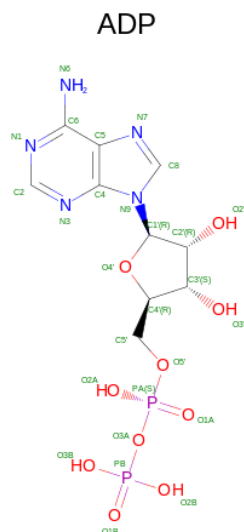
- Molecule 7 is a DNA chain called DNA (167-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	J	146	Total	C	N	O	P	0	0
			3011	1425	564	876	146		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

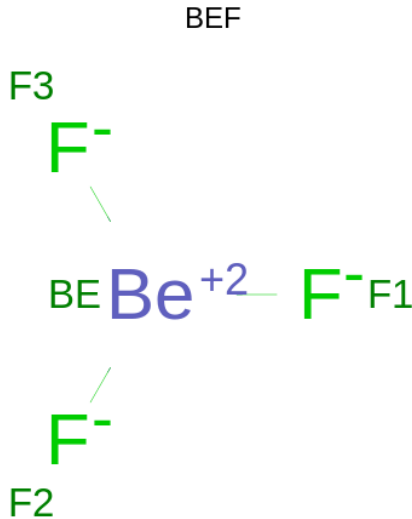
Mol	Chain	Residues	Atoms		AltConf
8	O	1	Total	Mg	0
			1	1	

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



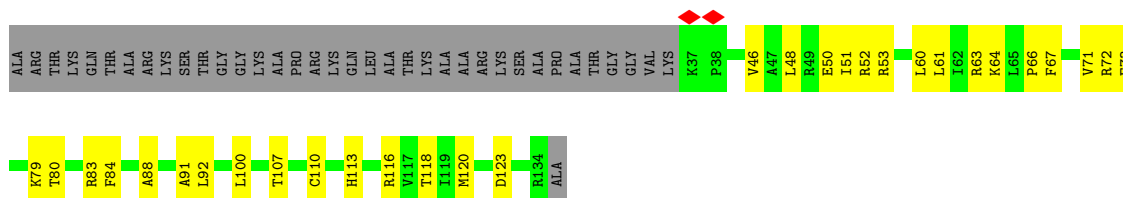
Mol	Chain	Residues	Atoms					AltConf
9	O	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 10 is BERYLLIUM TRIFLUORIDE ION (CCD ID: BEF) (formula: BeF_3).

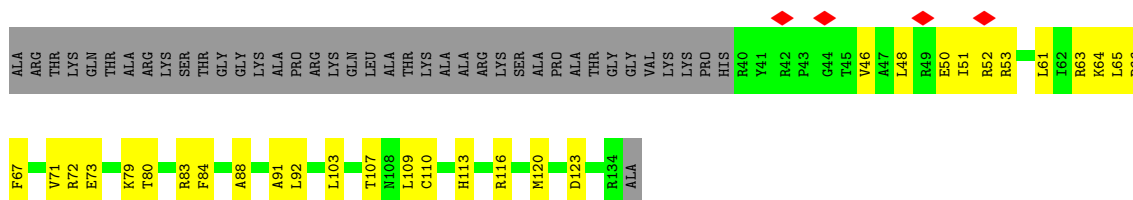


Mol	Chain	Residues	Atoms			AltConf
10	O	1	Total 4	Be 1	F 3	0

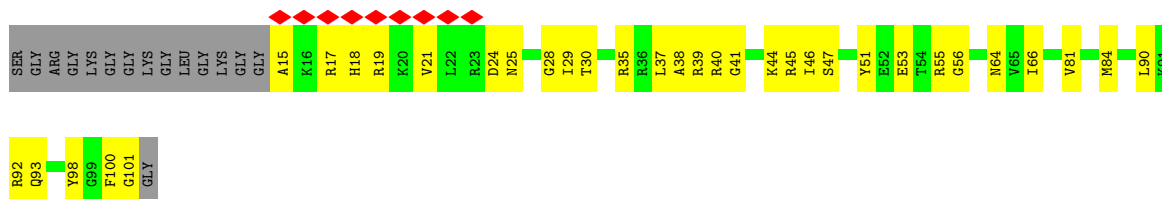
- Molecule 2: Histone H3.2



- Molecule 2: Histone H3.2



- Molecule 3: Histone H4

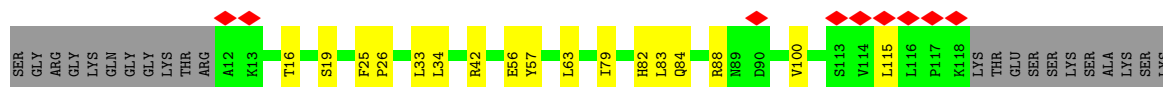


- Molecule 3: Histone H4

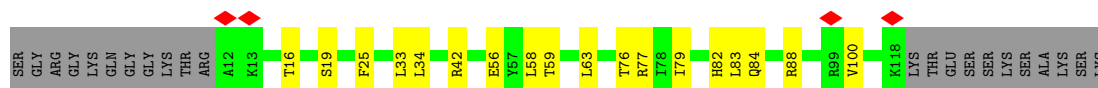




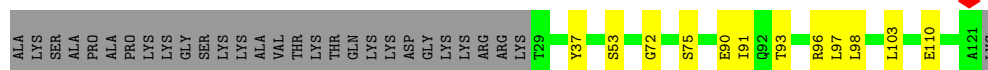
- Molecule 4: Histone H2A



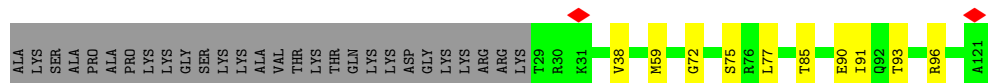
- Molecule 4: Histone H2A



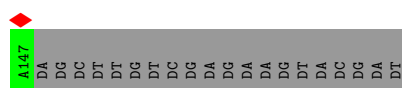
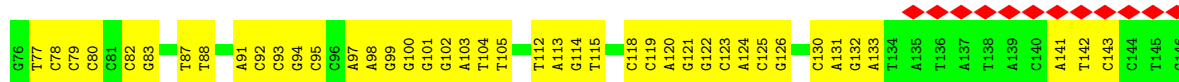
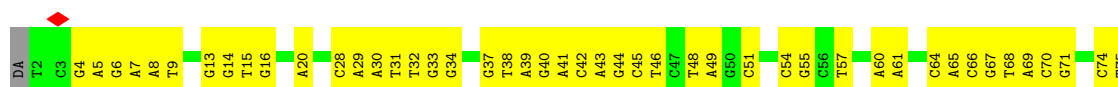
- Molecule 5: Histone H2B 1.1



- Molecule 5: Histone H2B 1.1



- Molecule 6: DNA (167-MER)



- Molecule 7: DNA (167-MER)

C127	T128	C129	G130	G131	C132	C134	A133	C135	G136	G137	T140	T141	C142	T143	C144	G145	A146	DT	DA	DT	DC	DC	DC	DC	DA	DA	DA	DC	DC	DC	DT	T1	C2	A3	G4	G5	A6	T7	G8	T9	A10	T11	A12	T13	A14	T15	C16	C20	A21	C22	G23	T24	G25	C26	C27	T28	G29	G30	A31	C34	T35	G39	A40	G41	T42	A43	A44	T45
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	82725	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 ($e^-/\text{\AA}^2$)	50.0	Depositor
Minimum defocus (nm)	1800	Depositor
Maximum defocus (nm)	3200	Depositor
Magnification	2250	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.314	Depositor
Minimum map value	-0.221	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.014	Depositor
Recommended contour level	0.0536	Depositor
Map size (Å)	211.20001, 211.20001, 211.20001	wwPDB
Map dimensions	160, 160, 160	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.32, 1.32, 1.32	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, MG, BEF

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	O	0.26	0/4799	0.54	1/6454 (0.0%)
2	A	0.58	0/813	0.68	0/1093
2	E	0.59	0/789	0.69	0/1059
3	B	0.68	0/711	0.80	0/950
3	F	0.66	0/680	0.72	0/912
4	C	0.57	0/821	0.59	0/1112
4	G	0.57	0/825	0.60	0/1116
5	D	0.61	0/729	0.66	0/985
5	H	0.60	0/737	0.66	0/993
6	I	0.49	0/3333	0.47	0/5137
7	J	0.48	0/3381	0.46	0/5221
All	All	0.48	0/17618	0.56	1/25032 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	O	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	959	PRO	N-CA-CB	7.26	110.24	103.41

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	O	810	TYR	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	O	4723	0	4808	316	0
2	A	801	0	831	28	0
2	E	779	0	815	27	0
3	B	703	0	757	29	0
3	F	672	0	698	20	0
4	C	811	0	849	16	0
4	G	815	0	860	16	0
5	D	718	0	725	10	0
5	H	726	0	747	9	0
6	I	2975	0	1639	79	0
7	J	3011	0	1639	90	0
8	O	1	0	0	0	0
9	O	27	0	12	8	0
10	O	4	0	0	4	0
All	All	16766	0	14380	592	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 19.

The worst 5 of 592 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:858:GLN:O	1:O:862:ARG:HB2	1.44	1.18
1:O:1253:LEU:O	1:O:1257:LEU:HB3	1.53	1.07
1:O:1343:ALA:O	1:O:1347:ASN:HB2	1.57	1.05
1:O:771:GLN:O	1:O:775:LEU:HB2	1.58	1.03
1:O:976:THR:O	1:O:980:ILE:HB	1.61	0.99

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	O	563/735 (77%)	481 (85%)	82 (15%)	0	100	100
2	A	96/135 (71%)	92 (96%)	4 (4%)	0	100	100
2	E	93/135 (69%)	89 (96%)	4 (4%)	0	100	100
3	B	85/102 (83%)	77 (91%)	8 (9%)	0	100	100
3	F	84/102 (82%)	82 (98%)	2 (2%)	0	100	100
4	C	105/129 (81%)	103 (98%)	2 (2%)	0	100	100
4	G	105/129 (81%)	103 (98%)	2 (2%)	0	100	100
5	D	91/122 (75%)	84 (92%)	7 (8%)	0	100	100
5	H	91/122 (75%)	84 (92%)	7 (8%)	0	100	100
All	All	1313/1711 (77%)	1195 (91%)	118 (9%)	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	O	520/667 (78%)	519 (100%)	1 (0%)	87	85
2	A	84/110 (76%)	84 (100%)	0	100	100
2	E	82/110 (74%)	82 (100%)	0	100	100
3	B	72/78 (92%)	72 (100%)	0	100	100
3	F	67/78 (86%)	67 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	C	81/101 (80%)	81 (100%)	0	100	100
4	G	82/101 (81%)	82 (100%)	0	100	100
5	D	77/102 (76%)	77 (100%)	0	100	100
5	H	79/102 (78%)	79 (100%)	0	100	100
All	All	1144/1449 (79%)	1143 (100%)	1 (0%)	87	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	1128	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	O	1189	GLN
3	B	25	ASN
4	G	38	ASN
1	O	1202	GLN
3	B	93	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
10	BEF	O	1503	-	0,3,3	-	-	-		
9	ADP	O	1502	-	27,29,29	1.32	4 (14%)	42,45,45	2.01	11 (26%)

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
9	ADP	O	1502	-	-	1/16/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	O	1502	ADP	C5-C4	4.30	1.47	1.39
9	O	1502	ADP	C5-C6	2.53	1.48	1.41
9	O	1502	ADP	C5-N7	-2.27	1.34	1.39
9	O	1502	ADP	C8-N7	2.21	1.35	1.31

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	O	1502	ADP	C5-C4-N3	-6.18	118.69	126.75
9	O	1502	ADP	N3-C4-N9	5.06	135.43	127.08
9	O	1502	ADP	C2-N3-C4	3.85	120.86	111.75
9	O	1502	ADP	PA-O3A-PB	-3.30	121.52	132.83
9	O	1502	ADP	N3-C2-N1	-3.18	123.62	128.60

There are no chirality outliers.

All (1) torsion outliers are listed below:

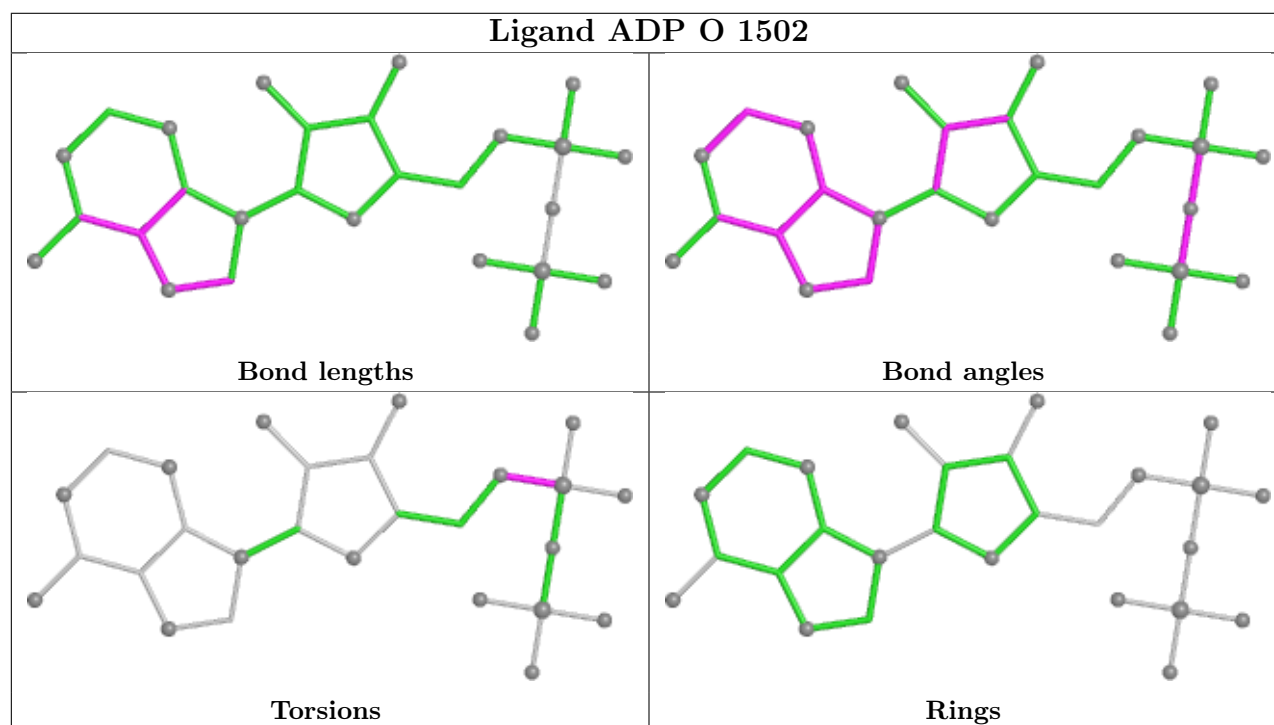
Mol	Chain	Res	Type	Atoms
9	O	1502	ADP	C5'-O5'-PA-O1A

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	O	1503	BEF	4	0
9	O	1502	ADP	8	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

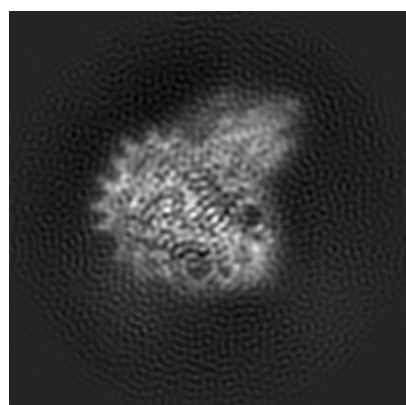
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6882. These allow visual inspection of the internal detail of the map and identification of artifacts.

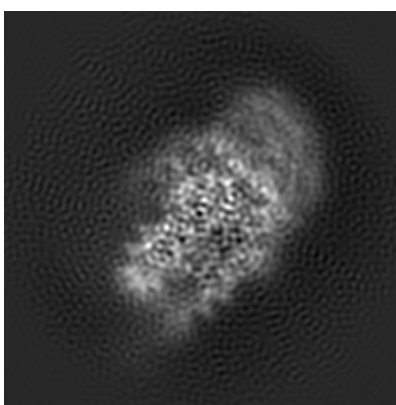
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

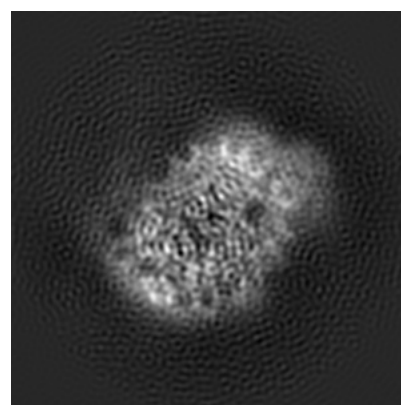
6.1.1 Primary map



X



Y

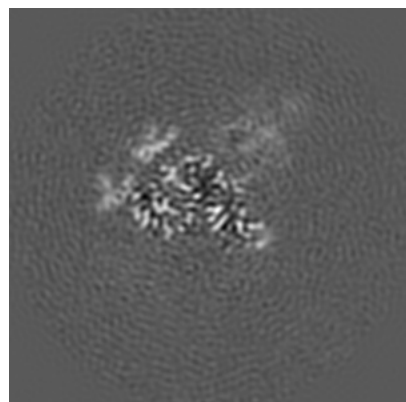


Z

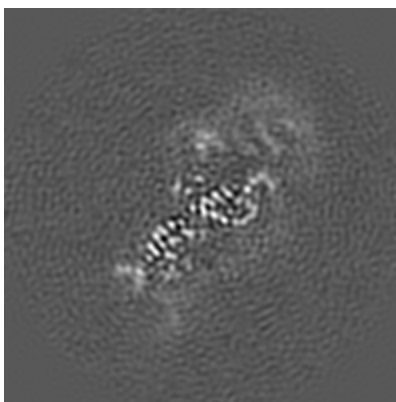
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

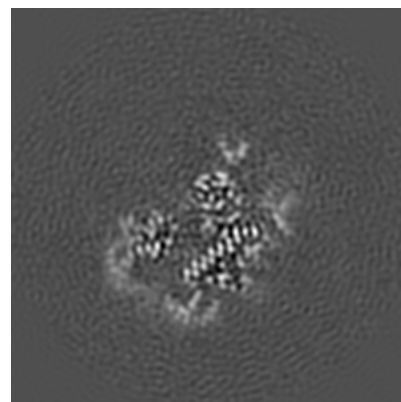
6.2.1 Primary map



X Index: 80



Y Index: 80

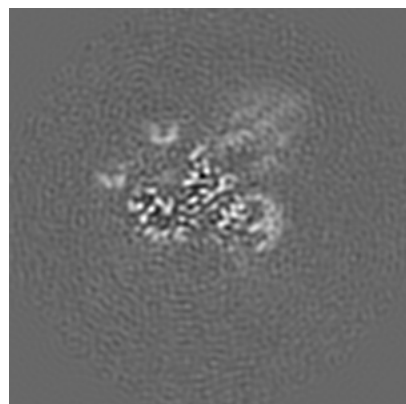


Z Index: 80

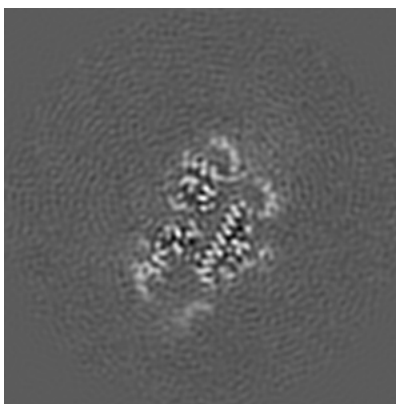
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

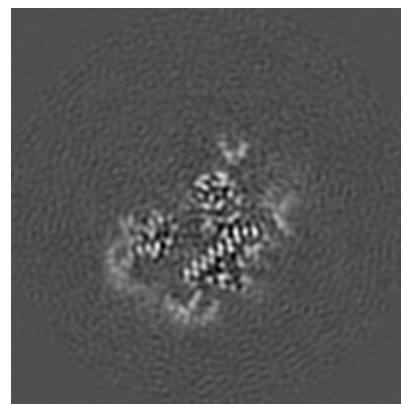
6.3.1 Primary map



X Index: 84



Y Index: 63

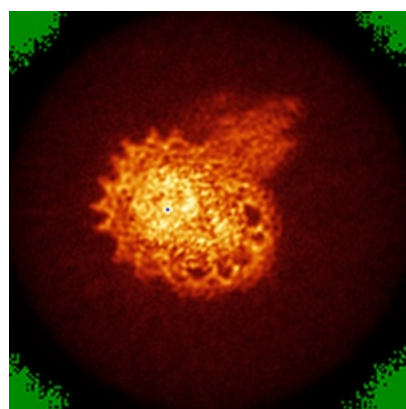


Z Index: 80

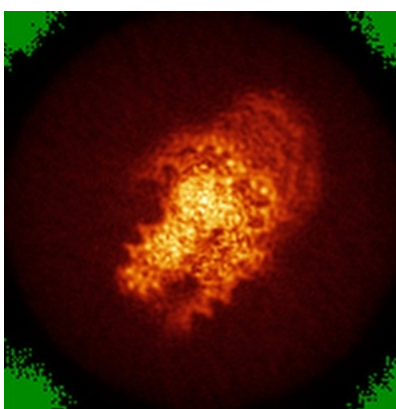
The images above show the largest variance slices of the map in three orthogonal directions.

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

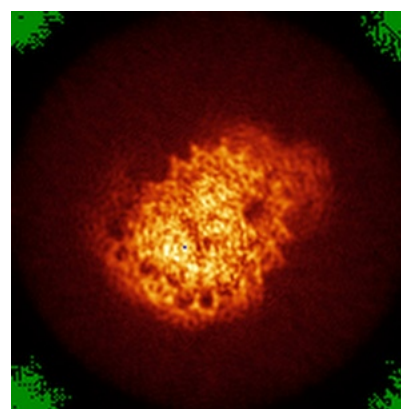
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0536. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

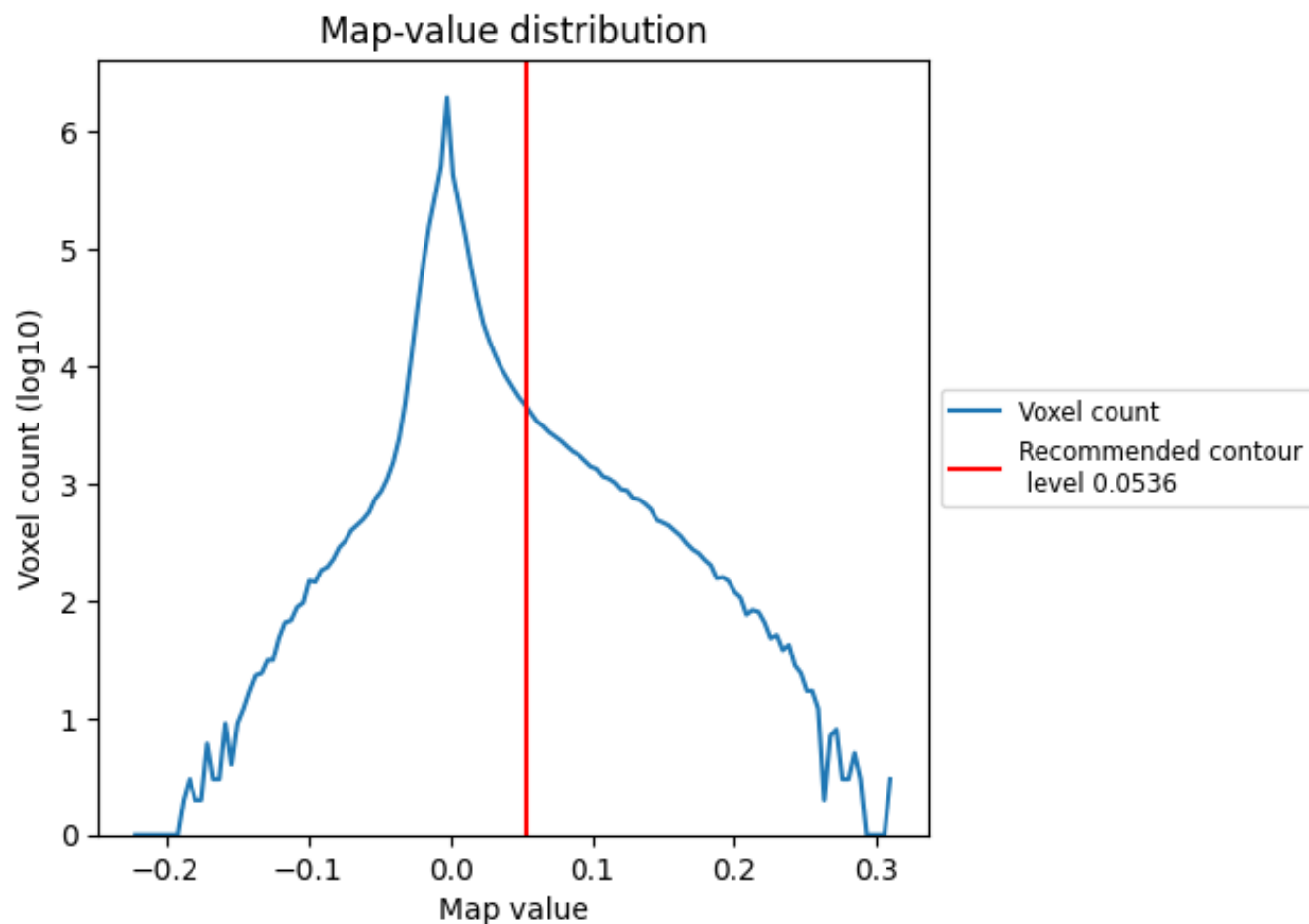
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

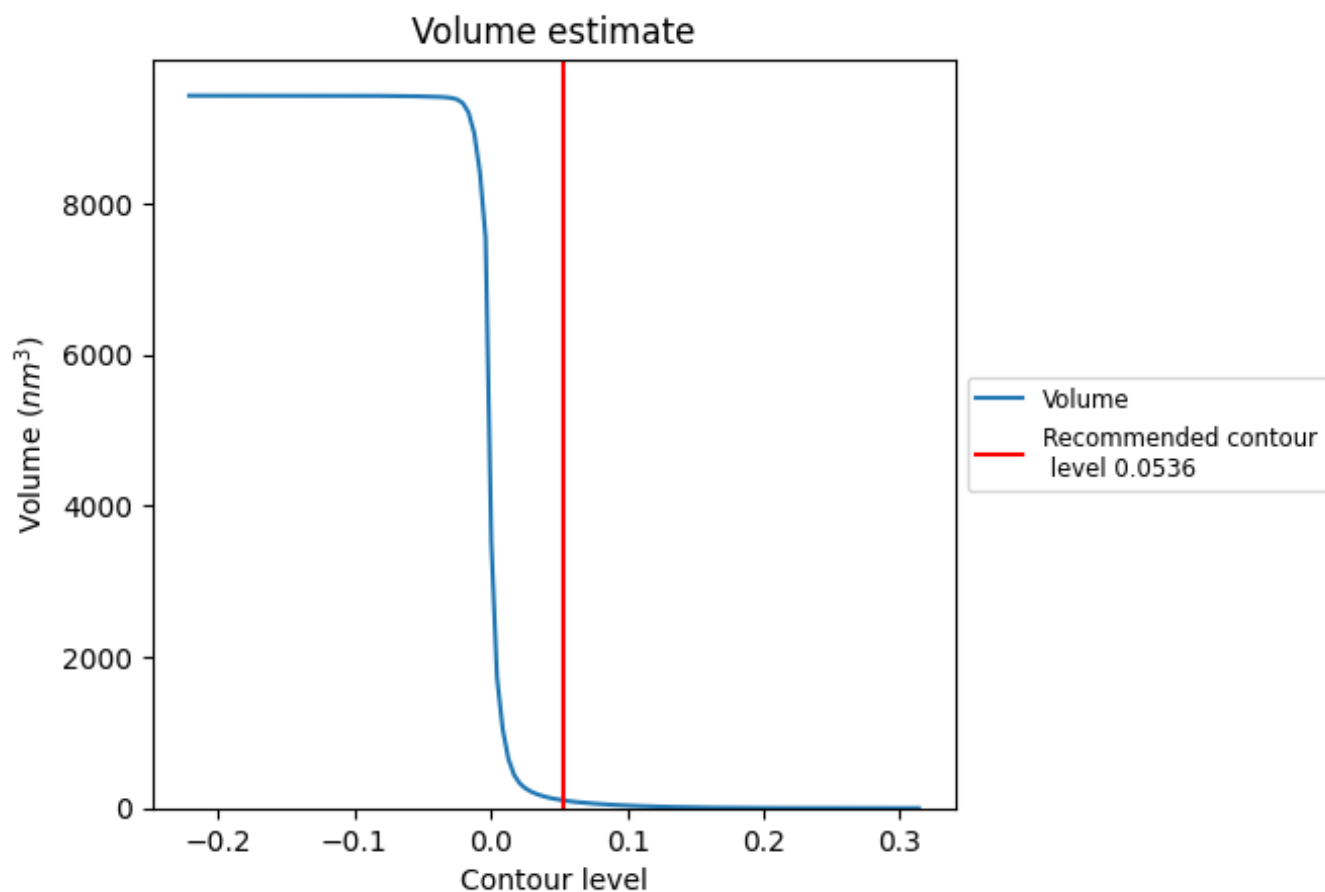
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

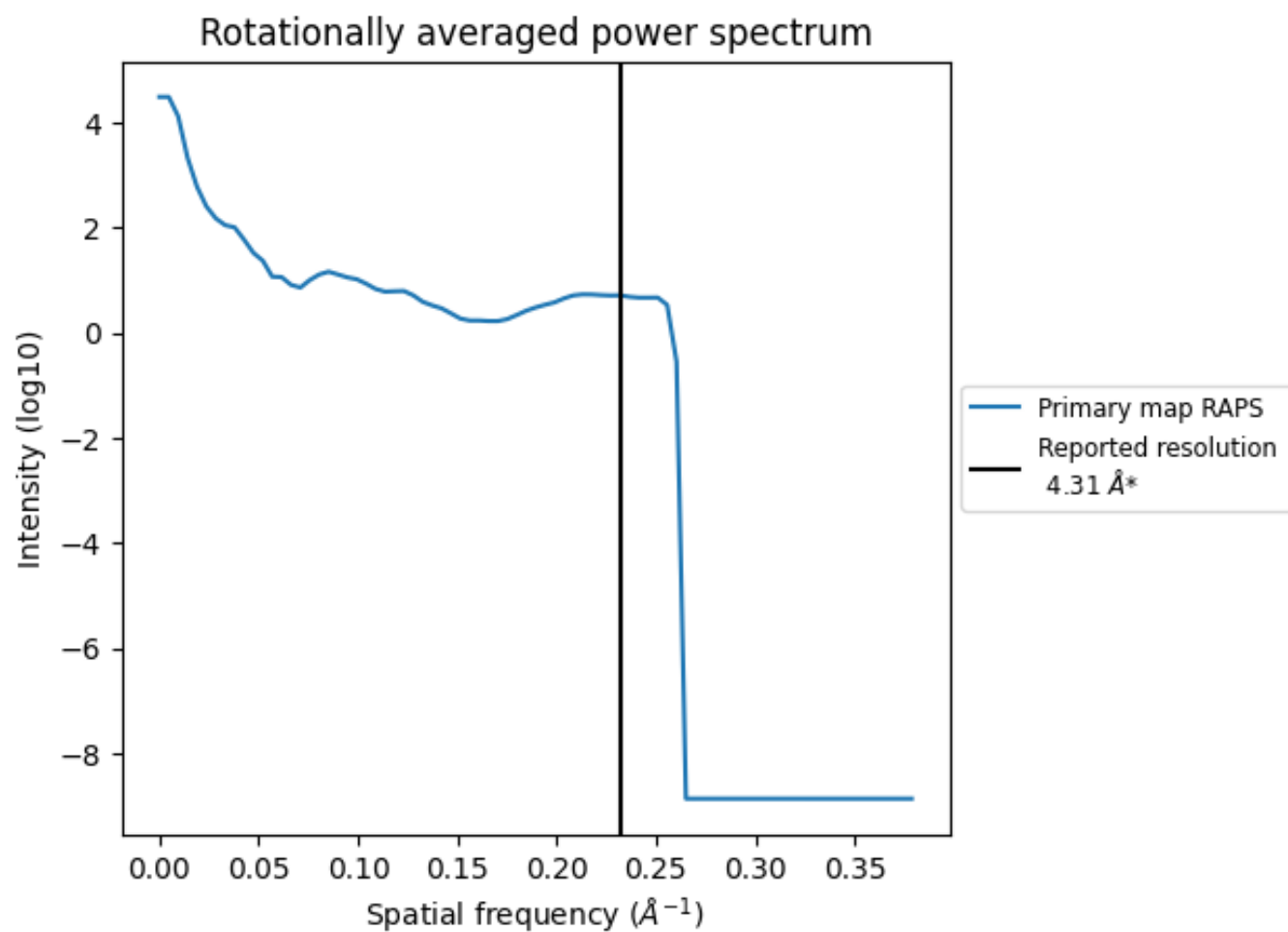
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 101 nm^3 ; this corresponds to an approximate mass of 91 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 [i](#)



*Reported resolution corresponds to spatial frequency of 0.232 Å⁻¹

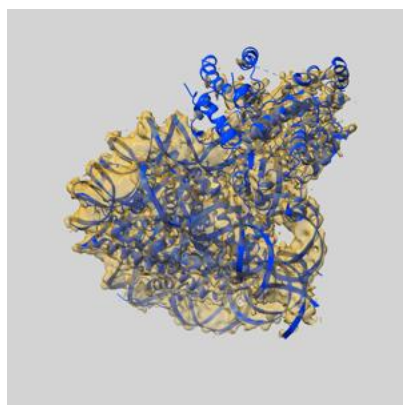
8 Fourier-Shell correlation ⓘ

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

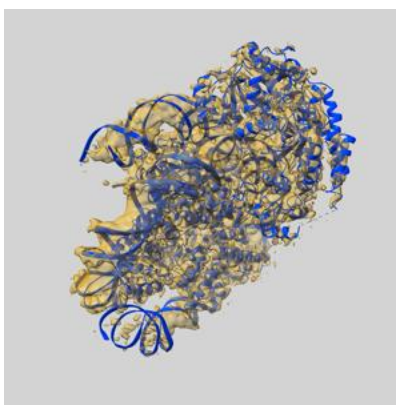
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-6882 and PDB model 5Z3U. Per-residue inclusion information can be found in section [3](#) on page [7](#).

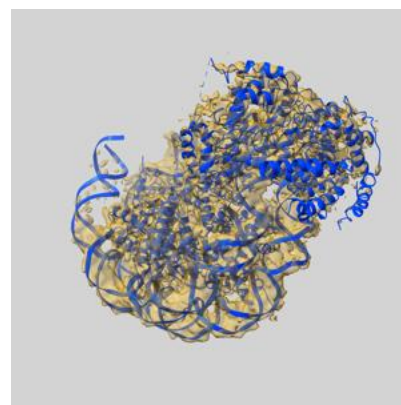
9.1 Map-model overlay [i](#)



X



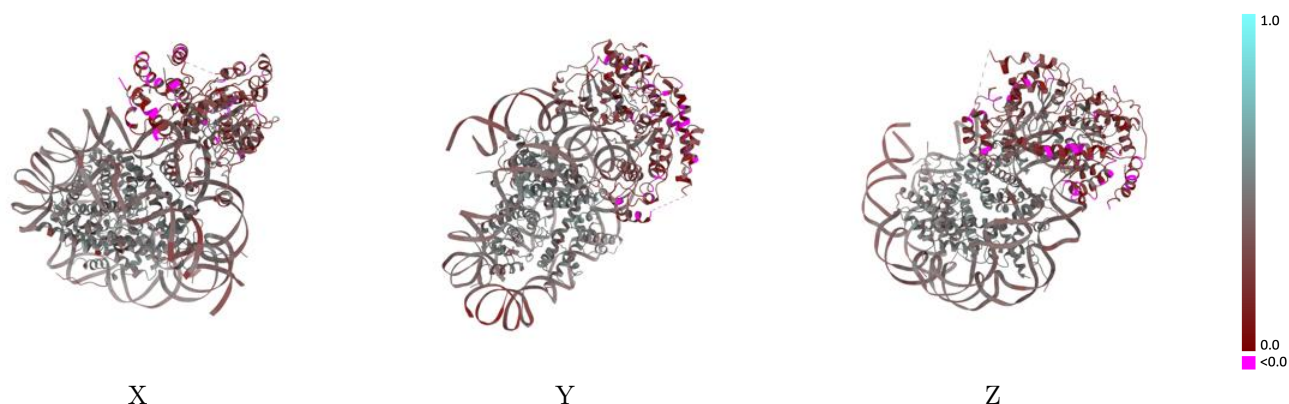
Y



Z

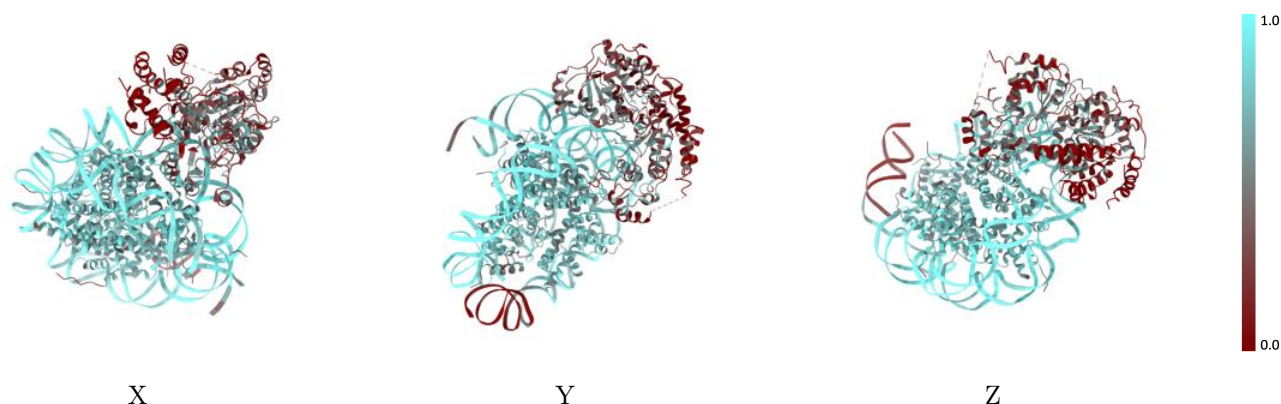
The images above show the 3D surface view of the map at the recommended contour level 0.0536 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



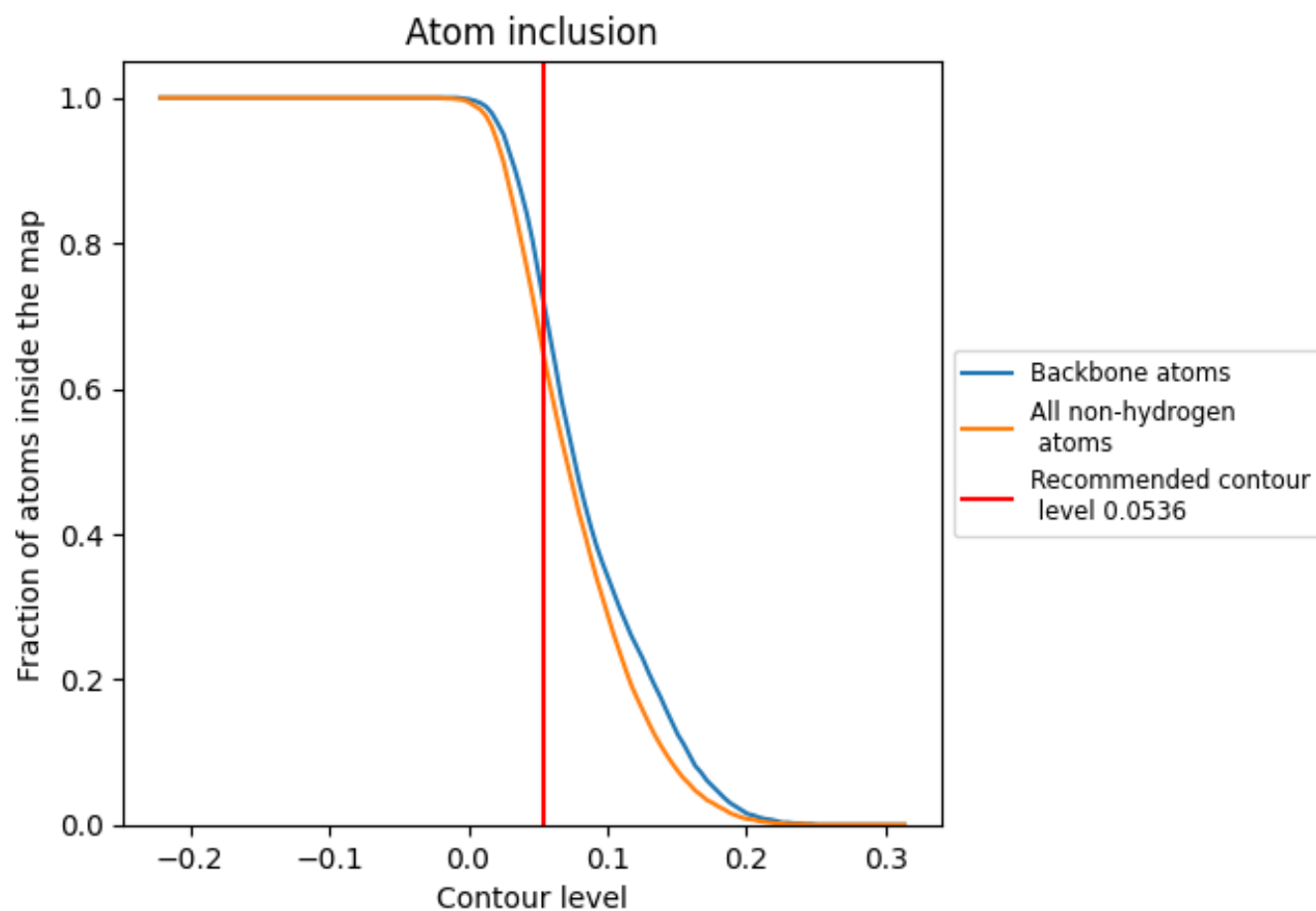
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0536).

9.4 Atom inclusion [i](#)



At the recommended contour level, 72% of all backbone atoms, 65% 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 (0.0536) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6510	0.3720
A	0.7560	0.4640
B	0.7340	0.4620
C	0.7410	0.4590
D	0.7790	0.4670
E	0.7560	0.4580
F	0.7560	0.4660
G	0.7770	0.4660
H	0.7810	0.4520
I	0.8210	0.3760
J	0.8160	0.3790
O	0.2980	0.2510

1.0

0.0

<0.0