

wwPDB EM Validation Summary Report (i)

Nov 20, 2022 – 03:11 pm GMT

PDB ID : 2WWB EMDB ID : EMD-1652

Title : CRYO-EM STRUCTURE OF THE MAMMALIAN SEC61 COMPLEX

BOUND TO THE ACTIVELY TRANSLATING WHEAT GERM 80S RI-

BOSOME

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Deposited on : 2009-10-22

Resolution : 6.48 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 $\frac{\text{https://www.wwpdb.org/validation/2017/EMValidationReportHelp}}{\text{with specific help available everywhere you see the (i) symbol.}$

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 (i)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev43

MolProbity: 4.02b-467

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $MapQ \quad : \quad 1.9.9$

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

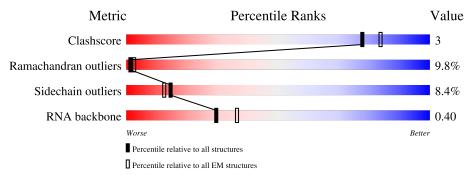
Validation Pipeline (wwPDB-VP) : 2.31.2

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 6.48 Å.

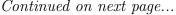
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive $(\# \mathrm{Entries})$	${ m EM\ structures} \ (\#{ m Entries})$
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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 <=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	Q	uality of chain	
			34%		
1	A	476		88%	9% ••
_	_		26%		
2	В	68		90%	9% •
		0.0	19%		
3	С	96	34% ••	63%	
	-		10%		
4	D	63	• 38%	51%	10%
_	-	2.4	<u>-</u>		_
5	Е	34	53%	4	4%
	-	25	8%		
6	F	25	72%		28%
_					
7	G	14	36%	64%	7





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Mol	Chain	Length		Quality of cha	ain	
8	Н	362	18%	16%	9% •	26%
9	I	184	15% 57%		19% 6%	• 17%
10	J	189	26% •		72%	
11	K	142	47%	9% •	42%	6
12	L	127	5%	74%		19% 6% •
13	M	113	11%		12% • •	26%
14	N	120	5% 44%	9% ••	42%	
15	О	51	24% 33%	31%		27%



2 Entry composition (i)

There are 15 unique types of molecules in this entry. The entry contains 14313 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 PROTEIN TRANSPORT PROTEIN SEC61 SUBUNIT ALPHA ISOFORM 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	476	Total 3675	C 2409	N 591	O 650	S 25	0	0

• Molecule 2 is a protein called PROTEIN TRANSPORT PROTEIN SEC61 SUBUNIT GAMMA.

Mol	Chain	Residues		Ato	ms	AltConf	Trace		
2	В	68	Total	С	N	О	S	0	0
			543	355	94	89	5		O

• Molecule 3 is a protein called PROTEIN TRANSPORT PROTEIN SEC61 SUBUNIT BETA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	36	Total	С	N	О	S	0	0
3		30	281	188	44	47	2	0	U

• Molecule 4 is a RNA chain called 5.8S RRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	63	Total 1347	C 603	N 245	O 436	P 63	0	0

• Molecule 5 is a RNA chain called 25S RRNA.

Mol	Chain	Residues		Atoms					Trace
5	E	34	Total 740	C 332	N 148	O 226	P 34	0	0

• Molecule 6 is a RNA chain called 25S RRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	25	Total 536	C 239	N 99	O 173	P 25	0	0



• Molecule 7 is a RNA chain called 25S RRNA.

Mol	Chain	Residues		Atoms Total C N O P 206 132 54 06 13				AltConf	Trace	
7	G	14	Total 296	C 133	N 54	O 96	P 13	0	0	

• Molecule 8 is a protein called 60S RIBOSOMAL PROTEIN L4-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Н	269	Total 2039	C 1281	N 391	O 363	S 4	0	0

• Molecule 9 is a protein called 60S RIBOSOMAL PROTEIN L17-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	153	Total 1212	C 756	N 236	O 219	S 1	0	0

• Molecule 10 is a protein called 60S RIBOSOMAL PROTEIN L19.

Mol	Chain	Residues		Ato	ms	AltConf	Trace		
10	J	53	Total 410	C 254	N 83	O 72	S 1	0	0

• Molecule 11 is a protein called 60S RIBOSOMAL PROTEIN L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	83	Total 663		N 111	O 126	S 2	0	0

• Molecule 12 is a protein called 60S RIBOSOMAL PROTEIN L26-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	127	Total 1002	C 630	N 193	O 178	S 1	0	0

• Molecule 13 is a protein called 60S RIBOSOMAL PROTEIN L31-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	84	Total 706	C 447	N 140	O 118	S 1	0	0

• Molecule 14 is a protein called 60S RIBOSOMAL PROTEIN L35.



Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	69	Total 547	C 345	N 101	O 99	S 2	0	0

 \bullet Molecule 15 is a protein called 60S RIBOSOMAL PROTEIN L39.

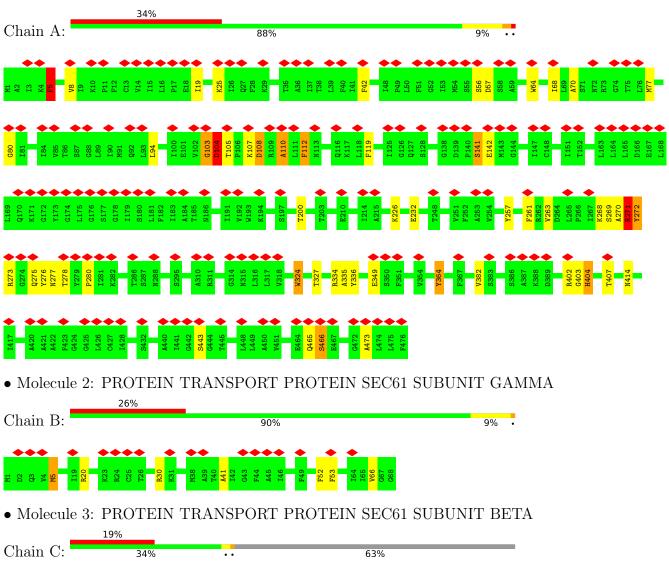
Mol	Chain	Residues		Ato	ms			AltConf	Trace
15	0	97	Total	С	N	О	S	0	0
15		31	316	200	66	48	2	0	U



3 Residue-property plots (i)

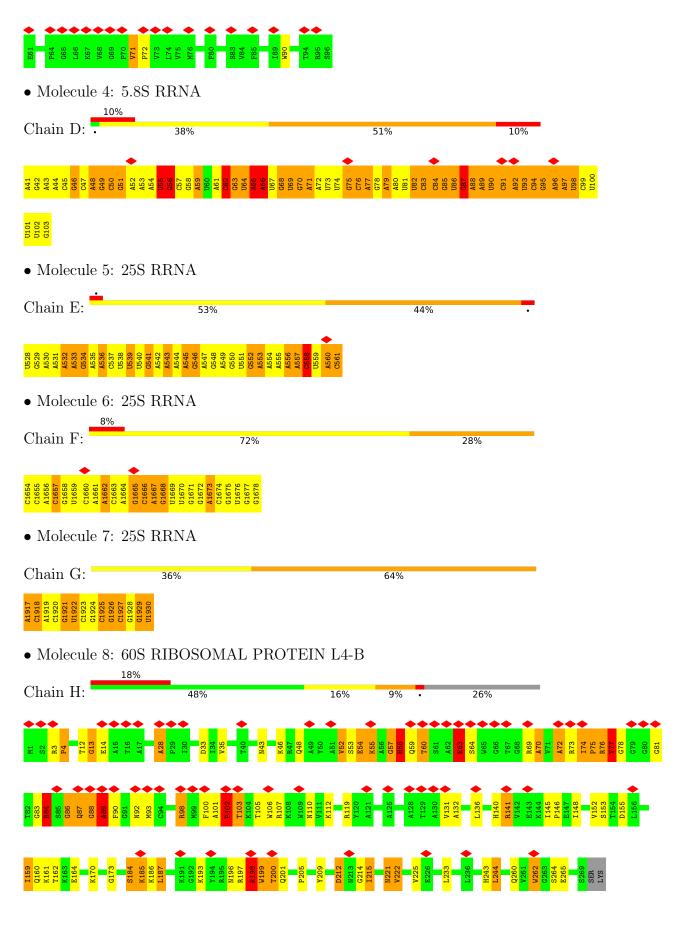
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: PROTEIN TRANSPORT PROTEIN SEC61 SUBUNIT ALPHA ISOFORM 1

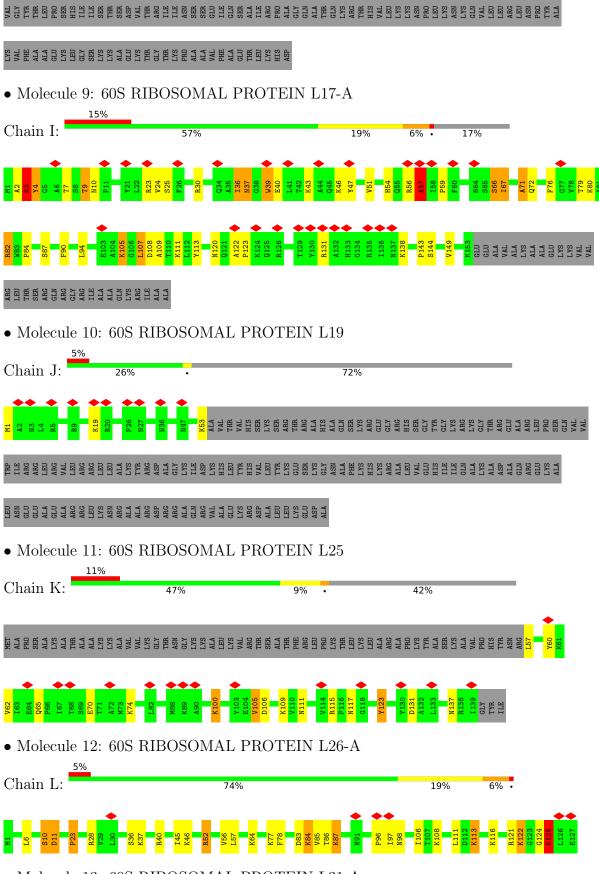




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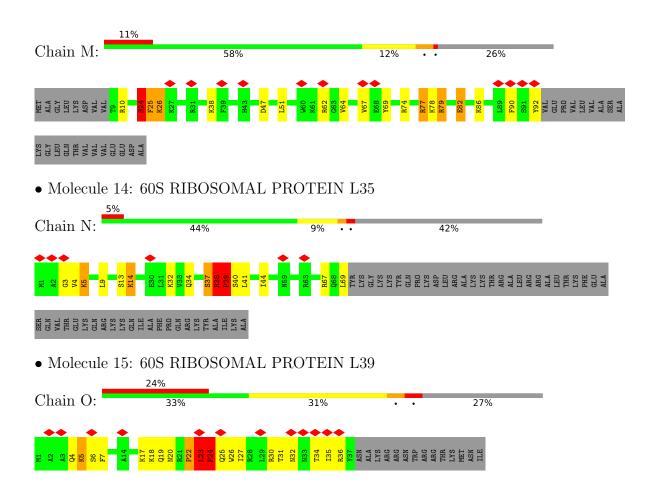






• Molecule 13: 60S RIBOSOMAL PROTEIN L31-A







4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	221445	Depositor
Resolution determination method	Not provided	
CTF correction method	DEFOCUS GROUP VOLUMES	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose $(e^-/\text{Å}^2)$	25	Depositor
Minimum defocus (nm)	1250	Depositor
Maximum defocus (nm)	4700	Depositor
Magnification	38000	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	1.706	Depositor
Minimum map value	-0.671	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.100	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	455.4, 455.4, 455.4	wwPDB
Map dimensions	368, 368, 368	wwPDB
Map angles (°)	90, 90, 90	wwPDB
Pixel spacing (Å)	1.2375, 1.2375, 1.2375	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	Bo	ond lengths	E	Bond angles
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	A	1.89	8/3755 (0.2%)	1.10	$22/5089 \ (0.4\%)$
2	В	0.94	0/553	1.12	2/738 (0.3%)
3	С	0.86	0/289	1.10	0/391
4	D	1.62	5/1508 (0.3%)	2.64	188/2348 (8.0%)
5	Е	1.67	2/833~(0.2%)	2.72	115/1298 (8.9%)
6	F	1.64	0/599	2.49	73/932 (7.8%)
7	G	1.70	2/330~(0.6%)	2.56	42/513 (8.2%)
8	Н	1.00	0/2079	1.46	33/2817 (1.2%)
9	I	1.03	0/1235	1.43	17/1662 (1.0%)
10	J	1.01	0/412	0.97	0/551
11	K	0.90	0/670	1.17	3/903 (0.3%)
12	L	1.02	0/1013	1.30	6/1351 (0.4%)
13	M	1.10	0/719	1.35	6/959~(0.6%)
14	N	1.05	1/549~(0.2%)	1.42	8/733 (1.1%)
15	O	1.10	0/321	1.45	4/426 (0.9%)
All	All	1.43	18/14865 (0.1%)	1.71	519/20711 (2.5%)

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
3	С	0	1
4	D	0	10
5	Е	0	6
6	F	0	3
7	G	0	1
8	Н	0	12
9	I	0	6
11	K	0	3
12	L	0	5

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Mol	Chain	#Chirality outliers	#Planarity outliers
13	M	0	2
14	N	0	2
15	O	0	7
All	All	0	63

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	Ideal(Å)
1	A	104	ASP	N-CA	81.65	3.09	1.46
1	A	5	PHE	CG-CD2	29.16	1.82	1.38
1	A	5	PHE	CG-CD1	28.74	1.81	1.38
1	A	5	PHE	CE1-CZ	22.75	1.80	1.37
1	A	5	PHE	CE2-CZ	22.22	1.79	1.37

The worst 5 of 519 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
5	Е	532	A	P-O3'-C3'	17.86	141.14	119.70
4	D	90	U	P-O3'-C3'	14.96	137.65	119.70
4	D	93	U	P-O3'-C3'	14.57	137.18	119.70
4	D	69	U	P-O3'-C3'	14.40	136.98	119.70
4	D	53	A	N1-C6-N6	14.32	127.19	118.60

There are no chirality outliers.

5 of 63 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	PHE	Sidechain
1	A	270	ALA	Peptide
1	A	271	ARG	Sidechain
1	A	336	TYR	Peptide
1	A	404	HIS	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	3675	0	3798	44	0
2	В	543	0	577	1	0
3	С	281	0	294	0	0
4	D	1347	0	678	2	0
5	Е	740	0	369	1	0
6	F	536	0	272	1	0
7	G	296	0	154	0	0
8	Н	2039	0	2106	16	0
9	I	1212	0	1231	8	0
10	J	410	0	452	0	0
11	K	663	0	699	4	0
12	L	1002	0	1093	2	0
13	M	706	0	741	1	0
14	N	547	0	613	8	0
15	О	316	0	349	2	0
All	All	14313	0	13426	81	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 3.

The worst 5 of 81 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:5:PHE:CD2	1:A:5:PHE:CE2	1.81	1.68
1:A:5:PHE:CD1	1:A:5:PHE:CE1	1.80	1.66
1:A:5:PHE:CE2	1:A:5:PHE:CZ	1.79	1.64
1:A:5:PHE:CD1	1:A:5:PHE:CG	1.81	1.61
1:A:5:PHE:CE1	1:A:5:PHE:CZ	1.80	1.61

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	474/476 (100%)	402 (85%)	46 (10%)	26 (6%)	2 19
2	В	$66/68 \; (97\%)$	60 (91%)	2 (3%)	4 (6%)	1 17
3	С	34/96~(35%)	27 (79%)	4 (12%)	3 (9%)	1 11
8	Н	267/362 (74%)	181 (68%)	38 (14%)	48 (18%)	0 2
9	I	151/184 (82%)	119 (79%)	15 (10%)	17 (11%)	0 7
10	J	51/189 (27%)	46 (90%)	4 (8%)	1 (2%)	7 38
11	K	81/142 (57%)	73 (90%)	5 (6%)	3 (4%)	3 24
12	L	125/127 (98%)	101 (81%)	9 (7%)	15 (12%)	0 6
13	M	82/113 (73%)	65 (79%)	8 (10%)	9 (11%)	0 7
14	N	67/120 (56%)	59 (88%)	3 (4%)	5 (8%)	1 13
15	О	35/51 (69%)	26 (74%)	0	9 (26%)	0 1
All	All	1433/1928 (74%)	1159 (81%)	134 (9%)	140 (10%)	1 9

5 of 140 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	SER
1	A	104	ASP
1	A	108	ASP
1	A	110	ALA
1	A	142	GLU

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	$398/398 \; (100\%)$	388 (98%)	10 (2%)	47 68
2	В	59/59~(100%)	57 (97%)	2 (3%)	37 60
3	С	32/74~(43%)	32 (100%)	0	100 100
8	Н	209/288~(73%)	179 (86%)	30 (14%)	3 16
9	I	124/146~(85%)	107 (86%)	17 (14%)	3 17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
10	J	44/154 (29%)	42 (96%)	2 (4%)	27 52
11	K	74/118 (63%)	67 (90%)	7 (10%)	8 28
12	L	110/110 (100%)	95 (86%)	15 (14%)	3 17
13	M	74/97 (76%)	64 (86%)	10 (14%)	4 18
14	N	62/105 (59%)	57 (92%)	5 (8%)	11 35
15	О	33/46 (72%)	28 (85%)	5 (15%)	3 15
All	All	1219/1595 (76%)	1116 (92%)	103 (8%)	14 34

5 of 103 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	I	144	SER
12	L	57	LEU
15	O	18	LYS
10	J	1	MET
11	K	111	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
8	Н	213	ASN
11	K	111	ASN
12	L	42	GLN
2	В	58	HIS
1	A	277	ASN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	D	62/63~(98%)	27 (43%)	10 (16%)
5	Е	33/34~(97%)	9 (27%)	3 (9%)
6	F	24/25~(96%)	3 (12%)	0
7	G	$14/14 \ (100\%)$	6 (42%)	2 (14%)
All	All	$133/136 \ (97\%)$	45 (33%)	15 (11%)

5 of 45 RNA backbone outliers are listed below:



Mol	Chain	Res	Type
4	D	48	A
4	D	49	G
4	D	51	G
4	D	56	G
4	D	59	A

5 of 15 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	D	90	U
7	G	1917	A
4	D	93	U
7	G	1926	G
5	Ε	558	G

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 monosaccharides 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.



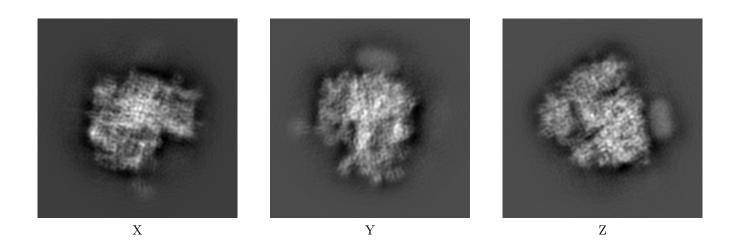
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-1652. 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)

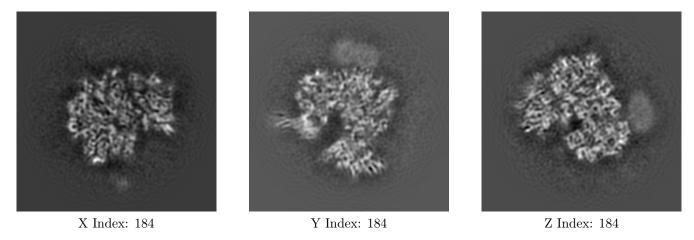
6.1.1 Primary map



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

6.2 Central slices (i)

6.2.1 Primary map

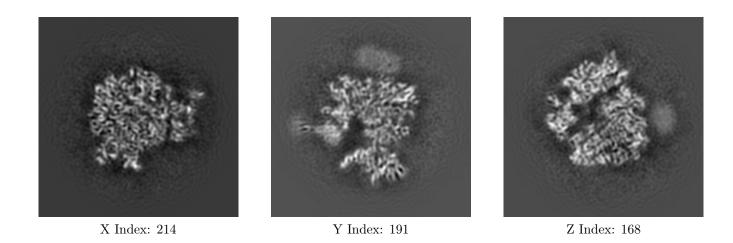




The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

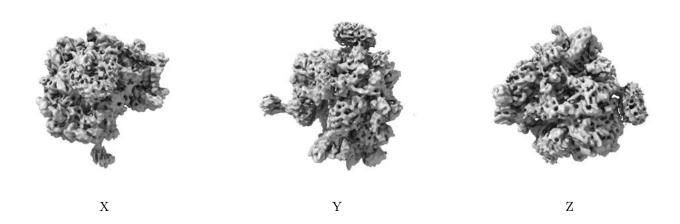
6.3.1 Primary map



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

6.4 Orthogonal surface views (i)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.



6.5 Mask visualisation (i)

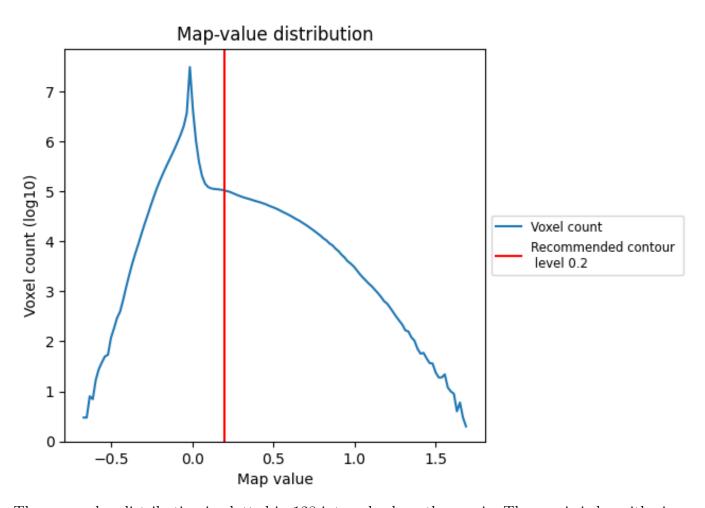
This section was not generated. No masks/segmentation were deposited.



7 Map analysis (i)

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

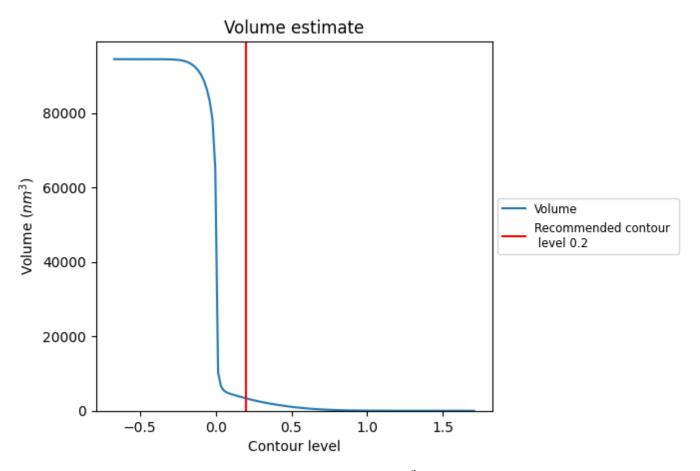
7.1 Map-value distribution (i)



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)

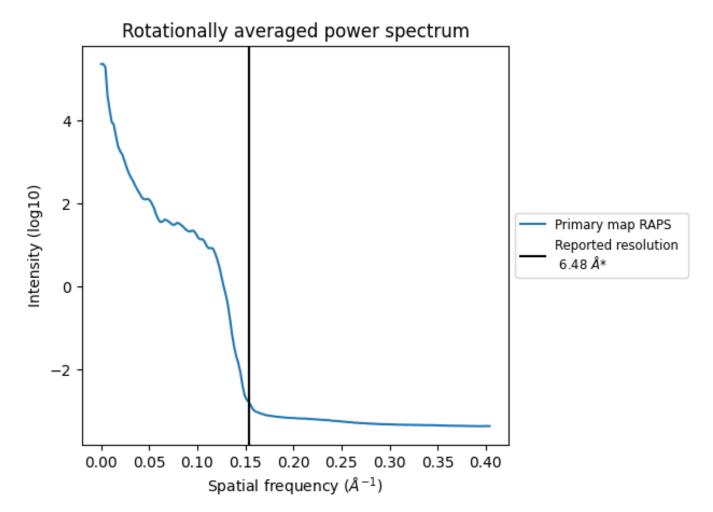


The volume at the recommended contour level is $3348~\mathrm{nm}^3$; this corresponds to an approximate mass of $3024~\mathrm{kDa}$.

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 0.154 $\rm \mathring{A}^{-1}$



8 Fourier-Shell correlation (i)

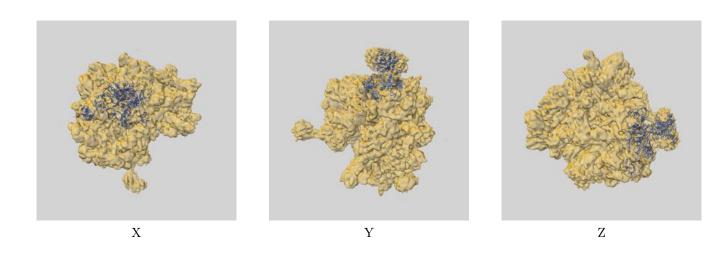
This section was not generated. No FSC curve or half-maps provided.



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-1652 and PDB model 2WWB. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

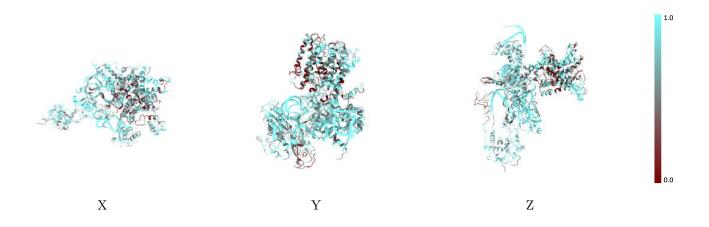


9.2 Q-score mapped to coordinate model (i)



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

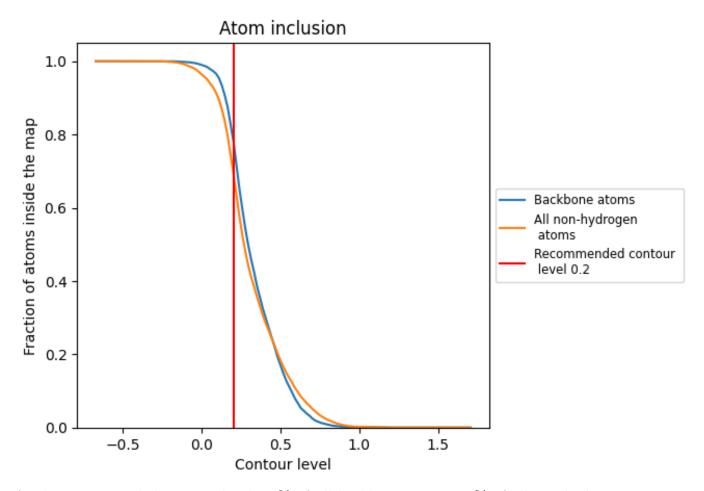
9.3 Atom inclusion mapped to coordinate model (i)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.2).



9.4 Atom inclusion (i)



At the recommended contour level, 79% of all backbone atoms, 70% of all non-hydrogen atoms, are inside the map.



9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6975	0.0920
A	0.5775	0.0860
В	0.6692	0.1140
С	0.3971	0.0410
D	0.8189	0.0880
E	0.9351	0.1460
F	0.8769	0.1340
G	0.9426	0.1730
Н	0.6628	0.0690
I	0.6888	0.0790
J	0.6667	0.1180
K	0.6503	0.0970
L	0.8564	0.0970
M	0.7074	0.0760
N	0.7425	0.0920
О	0.5691	0.0980



