



wwPDB EM Validation Summary Report ⓘ

Feb 25, 2024 – 10:29 PM EST

PDB ID : 6W6V
EMDB ID : EMD-21564
Title : Structure of yeast RNase MRP holoenzyme
Authors : Perederina, A.; Li, D.; Lee, H.; Bator, C.; Berezin, I.; Hafenstein, S.L.; Krasilnikov, A.S.
Deposited on : 2020-03-17
Resolution : 3.00 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:

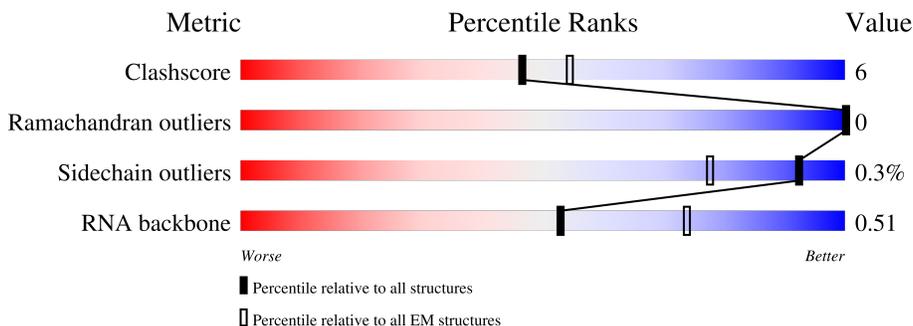
EMDB validation analysis : 0.0.1.dev70
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

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 3.00 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)
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 $\leq 5\%$. 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	340	
2	B	875	
3	D	279	
4	E	173	
5	F	158	
6	G	140	
7	H	133	

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Mol	Chain	Length	Quality of chain
8	I	293	
8	J	293	
9	K	198	
10	L	201	

2 Entry composition i

There are 10 unique types of molecules in this entry. The entry contains 24850 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 RNA chain called RNA component of RNase MRP NME1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	A	294	6261	2805	1097	2064	295	1	0

- Molecule 2 is a protein called Ribonucleases P/MRP protein subunit POP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	780	6345	4015	1154	1140	36	0	0

- Molecule 3 is a protein called RNases MRP/P 32.9 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	D	212	1759	1131	300	320	8	0	0

- Molecule 4 is a protein called Ribonuclease P/MRP protein subunit POP5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	E	169	1335	837	237	253	8	0	0

- Molecule 5 is a protein called Ribonucleases P/MRP protein subunit POP6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	F	158	1280	809	223	243	5	0	0

- Molecule 6 is a protein called Ribonucleases P/MRP protein subunit POP7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	G	126	1006	635	177	192	2	0	0

- Molecule 7 is a protein called Ribonucleases P/MRP protein subunit POP8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	H	125	1033	655	169	207	2	0	0

- Molecule 8 is a protein called Ribonuclease P/MRP protein subunit RPP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	I	243	1891	1194	342	346	9	0	0
8	J	293	2260	1415	413	422	10	0	0

- Molecule 9 is a protein called Ribonuclease MRP protein subunit SNM1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	K	79	648	411	117	116	4	0	0

- Molecule 10 is a protein called Ribonuclease MRP protein subunit RMP1.

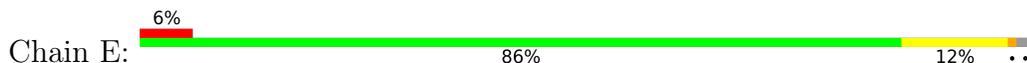
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	L	121	1032	665	189	172	6	0	0



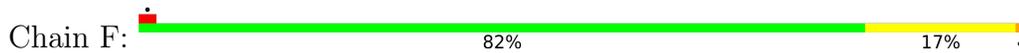
• Molecule 3: RNases MRP/P 32.9 kDa subunit



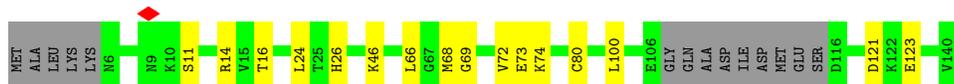
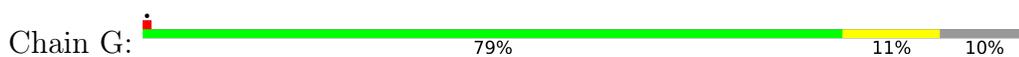
• Molecule 4: Ribonuclease P/MRP protein subunit POP5



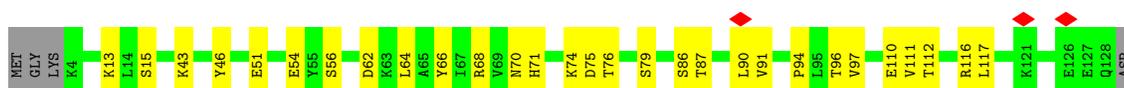
• Molecule 5: Ribonucleases P/MRP protein subunit POP6



• Molecule 6: Ribonucleases P/MRP protein subunit POP7



• Molecule 7: Ribonucleases P/MRP protein subunit POP8



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	155205	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 ($e^-/\text{\AA}^2$)	39	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	3.131	Depositor
Minimum map value	-1.597	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.094	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	340.48, 340.48, 340.48	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.665, 0.665, 0.665	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	0.35	0/6993	0.76	0/10870
2	B	0.29	0/6483	0.47	1/8734 (0.0%)
3	D	0.28	0/1790	0.47	0/2401
4	E	0.28	0/1351	0.45	0/1814
5	F	0.27	0/1298	0.48	0/1743
6	G	0.28	0/1019	0.55	1/1368 (0.1%)
7	H	0.28	0/1055	0.47	0/1428
8	I	0.27	0/1919	0.49	0/2592
8	J	0.26	0/2291	0.46	0/3092
9	K	0.27	0/662	0.53	0/891
10	L	0.28	0/1053	0.43	0/1418
All	All	0.30	0/25914	0.58	2/36351 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
6	G	66	LEU	CA-CB-CG	6.26	129.70	115.30
2	B	626	ARG	CA-CB-CG	5.15	124.73	113.40

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	6261	0	3165	59	0
2	B	6345	0	6427	94	0
3	D	1759	0	1817	21	0
4	E	1335	0	1380	21	0
5	F	1280	0	1317	17	0
6	G	1006	0	1048	14	0
7	H	1033	0	986	22	0
8	I	1891	0	1982	30	0
8	J	2260	0	2351	41	0
9	K	648	0	649	11	0
10	L	1032	0	1048	17	0
All	All	24850	0	22170	304	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 6.

The worst 5 of 304 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:A:167:G:H1	1:A:176:U:H3	1.28	0.80
5:F:47:HIS:HB3	5:F:81:ILE:HG22	1.65	0.79
1:A:79:G:H22	4:E:3:ARG:HH22	1.34	0.75
1:A:164:G:H1	1:A:179:U:H3	1.35	0.75
1:A:303:C:OP2	8:J:187:ARG:NH2	2.22	0.73

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	
2	B	772/875 (88%)	690 (89%)	82 (11%)	0	100	100
3	D	210/279 (75%)	189 (90%)	21 (10%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	E	167/173 (96%)	146 (87%)	21 (13%)	0	100	100
5	F	156/158 (99%)	142 (91%)	14 (9%)	0	100	100
6	G	122/140 (87%)	113 (93%)	9 (7%)	0	100	100
7	H	123/133 (92%)	105 (85%)	18 (15%)	0	100	100
8	I	241/293 (82%)	211 (88%)	30 (12%)	0	100	100
8	J	291/293 (99%)	259 (89%)	32 (11%)	0	100	100
9	K	77/198 (39%)	64 (83%)	13 (17%)	0	100	100
10	L	117/201 (58%)	108 (92%)	9 (8%)	0	100	100
All	All	2276/2743 (83%)	2027 (89%)	249 (11%)	0	100	100

There are no Ramachandran outliers to report.

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	
2	B	704/785 (90%)	703 (100%)	1 (0%)	93	98
3	D	196/261 (75%)	195 (100%)	1 (0%)	88	96
4	E	156/160 (98%)	155 (99%)	1 (1%)	86	95
5	F	149/149 (100%)	147 (99%)	2 (1%)	69	89
6	G	116/127 (91%)	116 (100%)	0	100	100
7	H	116/123 (94%)	116 (100%)	0	100	100
8	I	217/258 (84%)	217 (100%)	0	100	100
8	J	258/258 (100%)	256 (99%)	2 (1%)	81	93
9	K	73/184 (40%)	73 (100%)	0	100	100
10	L	110/182 (60%)	110 (100%)	0	100	100
All	All	2095/2487 (84%)	2088 (100%)	7 (0%)	92	97

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	F	57	LYS
5	F	132	ARG
8	J	144	LYS
8	J	52	LYS
4	E	168	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
5	F	111	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	285/340 (83%)	67 (23%)	0

5 of 67 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	16	A
1	A	26	A
1	A	35	U
1	A	36	U
1	A	39	A

There are no RNA pucker outliers to report.

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.

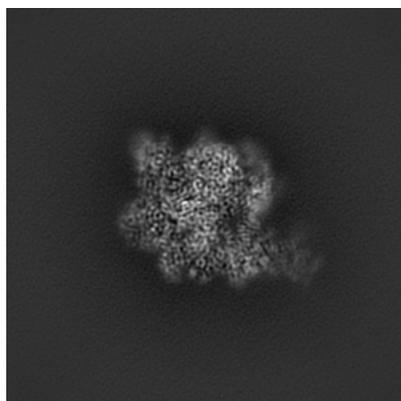
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-21564. These allow visual inspection of the internal detail of the map and identification of artifacts.

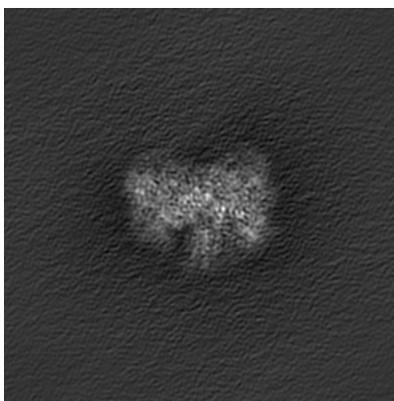
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

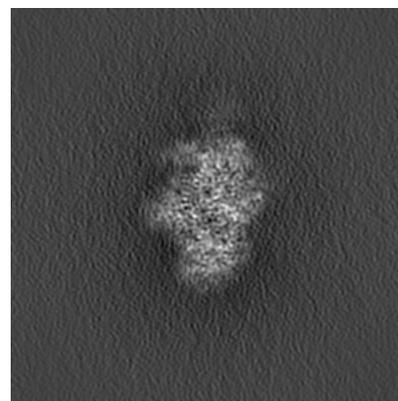
6.1.1 Primary map



X



Y

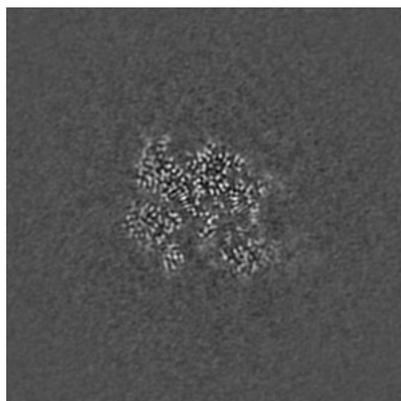


Z

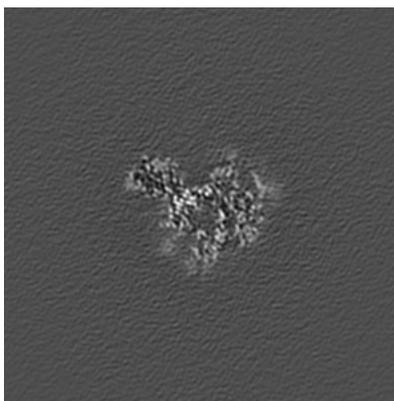
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

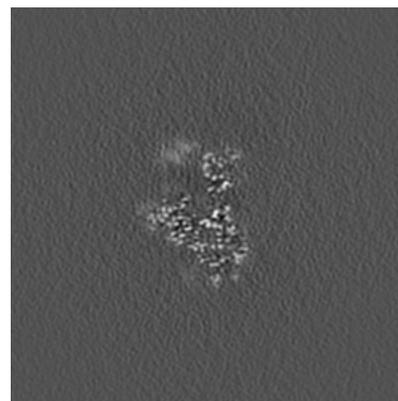
6.2.1 Primary map



X Index: 256



Y Index: 256

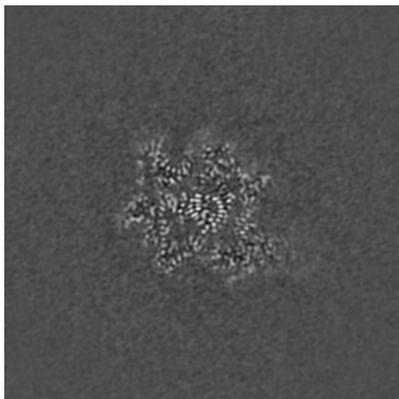


Z Index: 256

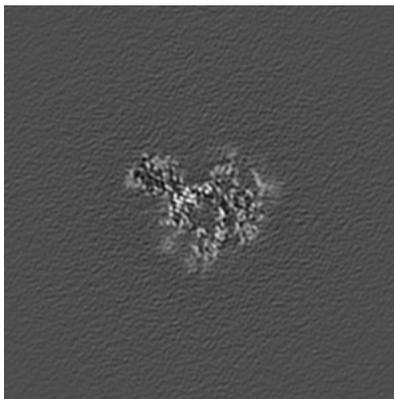
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

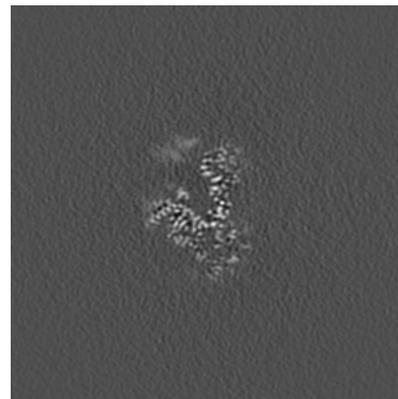
6.3.1 Primary map



X Index: 267



Y Index: 256

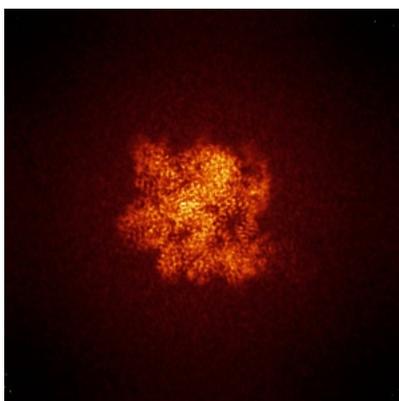


Z Index: 260

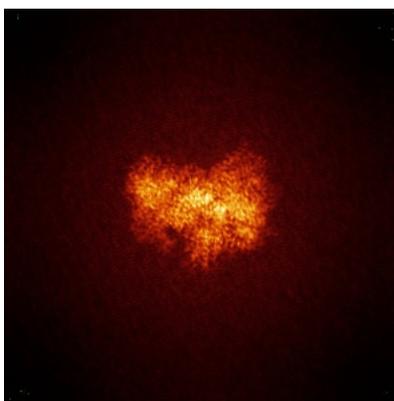
The images above show the largest variance slices of the map in three orthogonal directions.

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

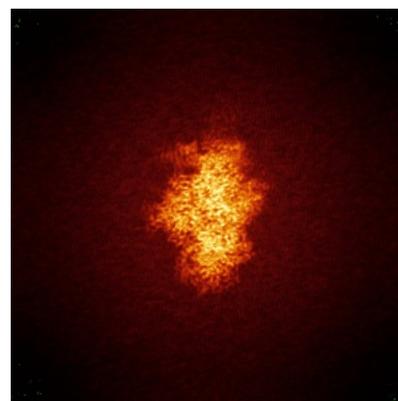
6.4.1 Primary map



X



Y

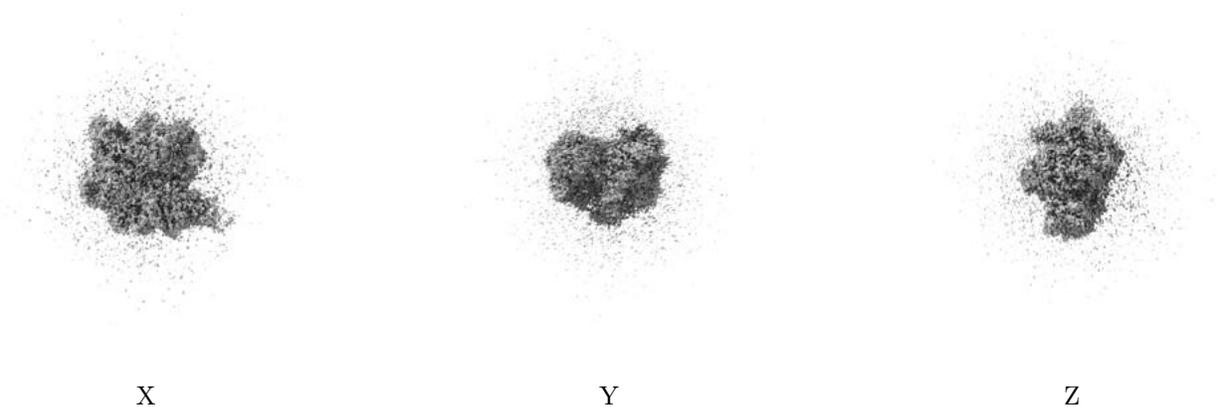


Z

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 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

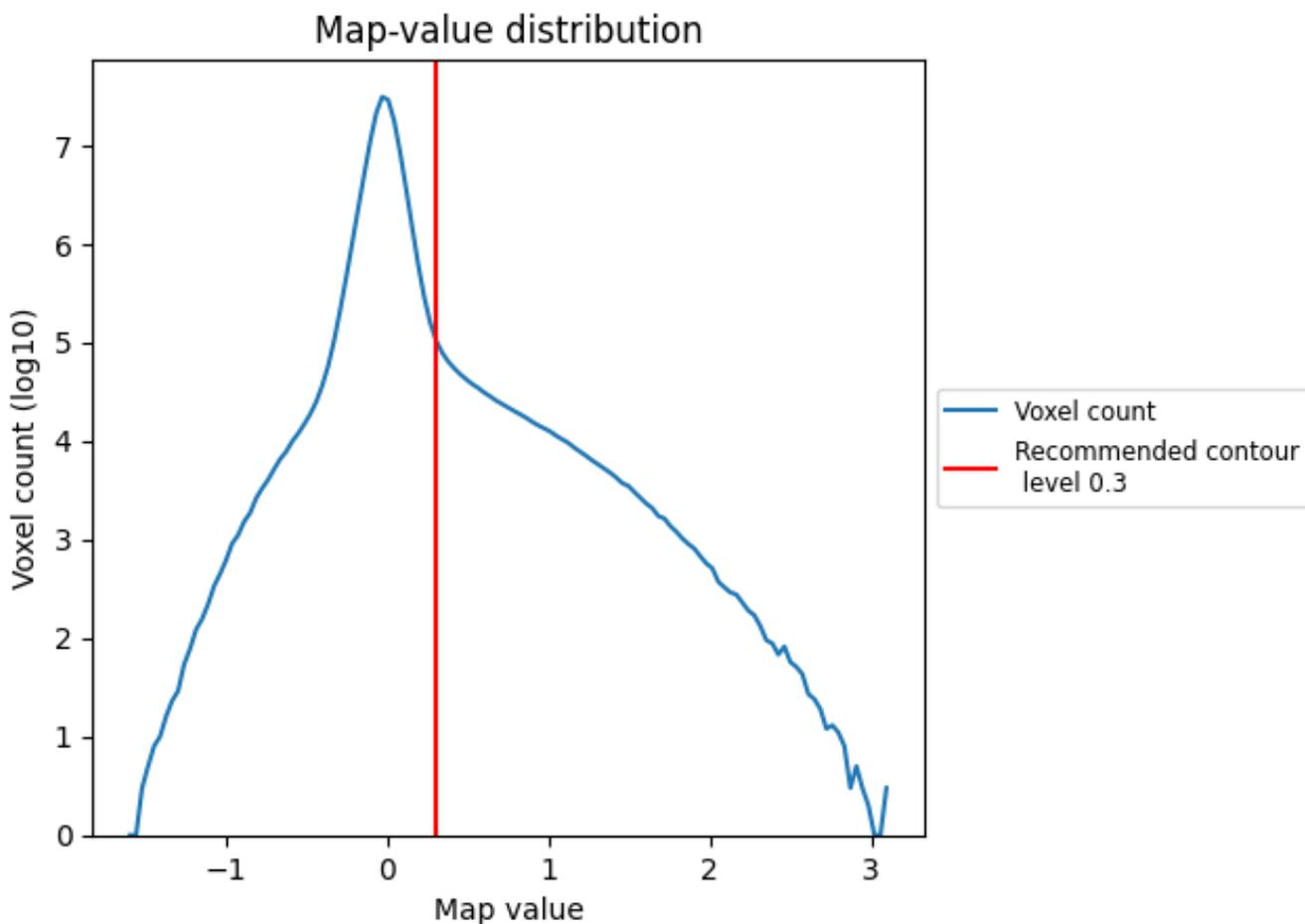
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

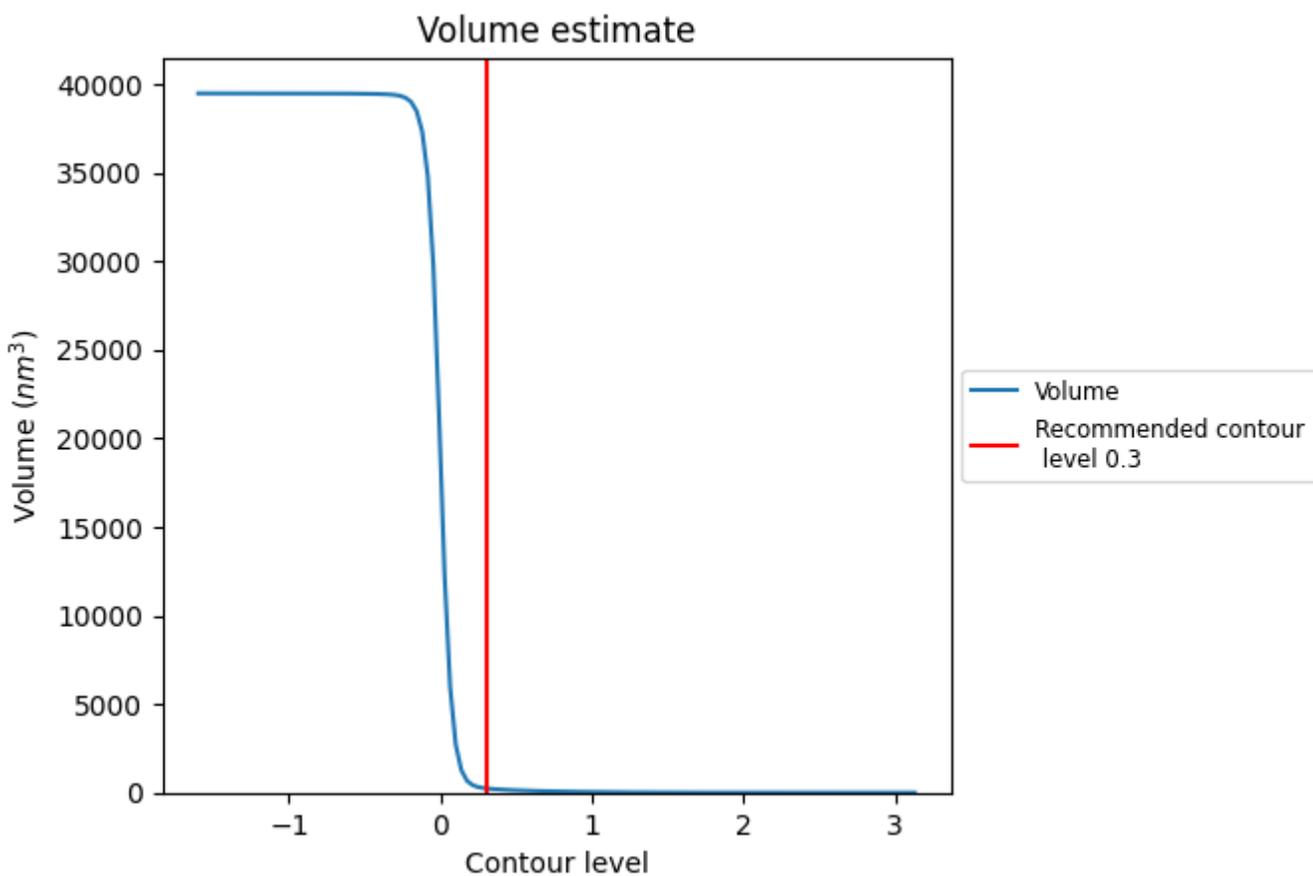
This section contains the results of statistical analysis of the map.

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.

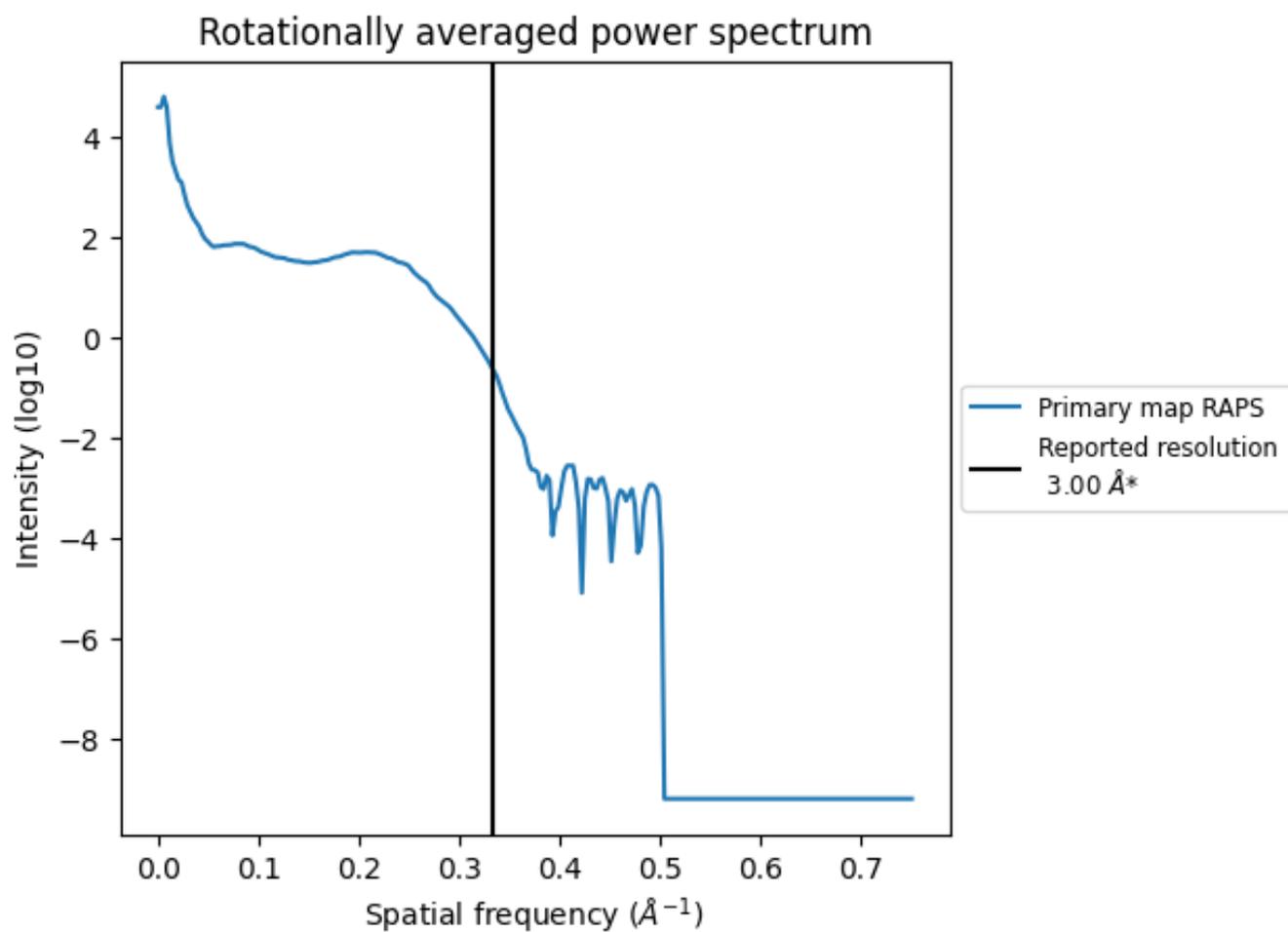
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 243 nm³; this corresponds to an approximate mass of 220 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 0.333 \AA^{-1}

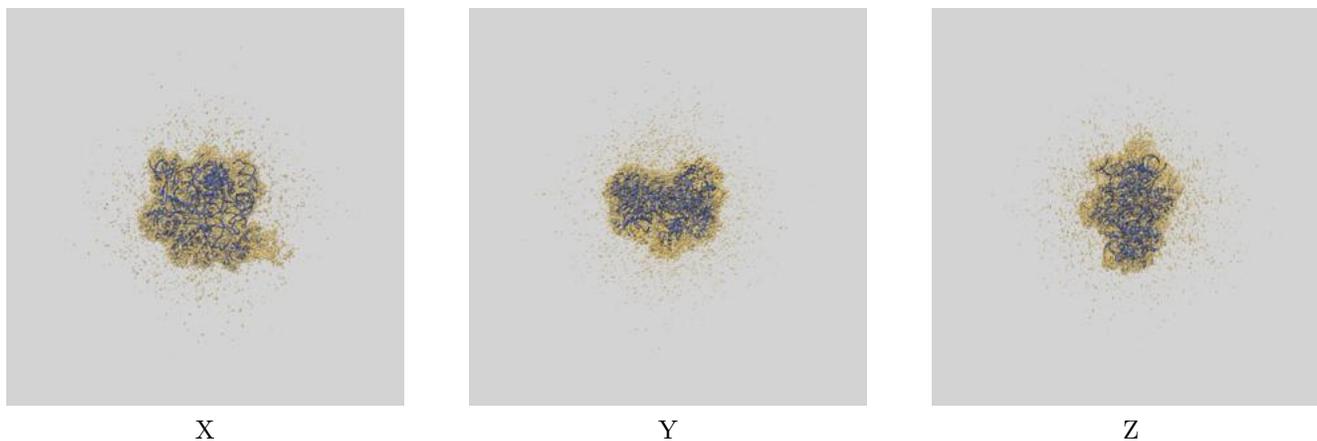
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

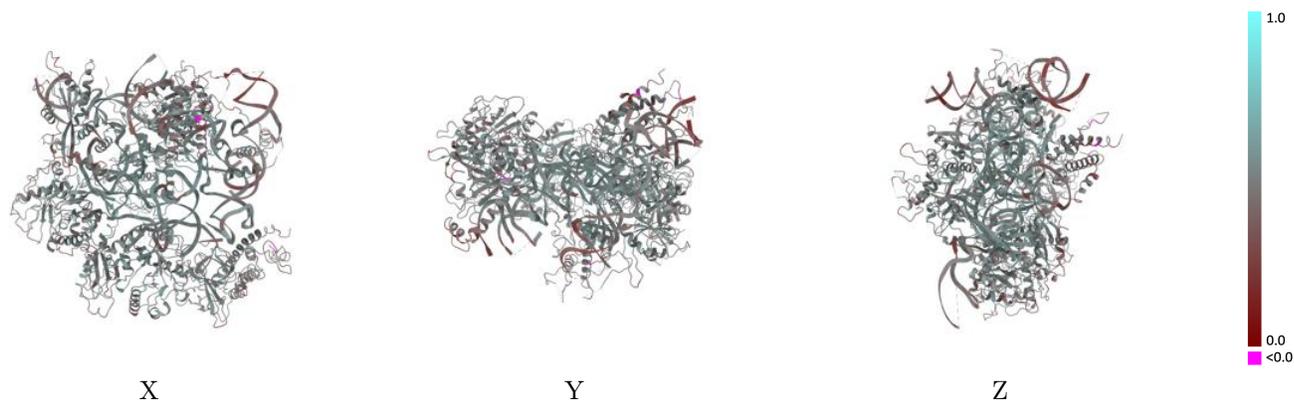
This section contains information regarding the fit between EMDB map EMD-21564 and PDB model 6W6V. Per-residue inclusion information can be found in section [3](#) on page [6](#).

9.1 Map-model overlay [i](#)



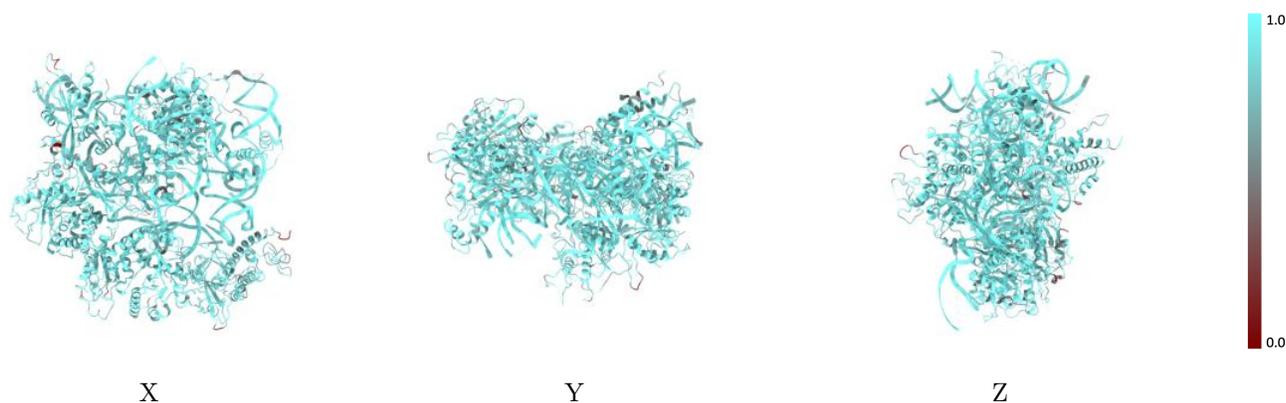
The images above show the 3D surface view of the map at the recommended contour level 0.3 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



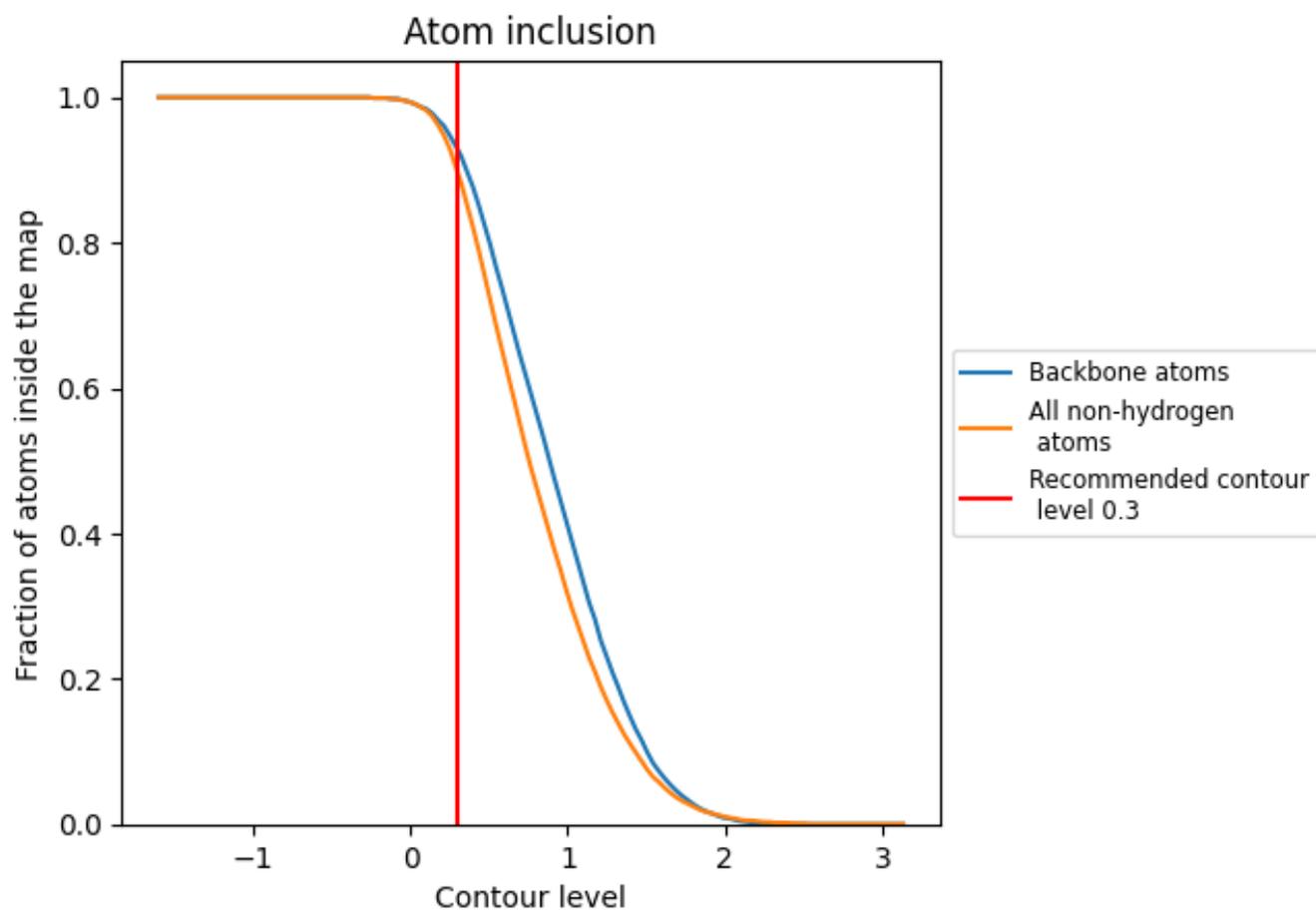
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.3).

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 90% 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 (0.3) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8980	 0.4890
A	 0.9350	 0.4840
B	 0.9070	 0.5060
D	 0.8510	 0.4810
E	 0.8500	 0.4830
F	 0.8910	 0.4810
G	 0.8940	 0.5110
H	 0.8600	 0.4600
I	 0.8970	 0.4910
J	 0.8950	 0.5010
K	 0.7740	 0.4470
L	 0.8830	 0.4590

