



wwPDB EM Validation Summary Report ⓘ

Mar 25, 2026 – 03:43 PM UTC

PDB ID : 5L1D / pdb_00005l1d
EMDB ID : EMD-8303
Title : Structure of rabbit RyR2 in complex with FKBP12.6 in a closed state (conformation C1)
Authors : Dhindwal, S.; Lobo, J.J.; Samso, M.
Deposited on : 2016-07-29
Resolution : 11.00 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:

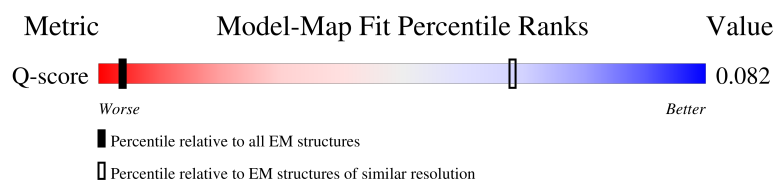
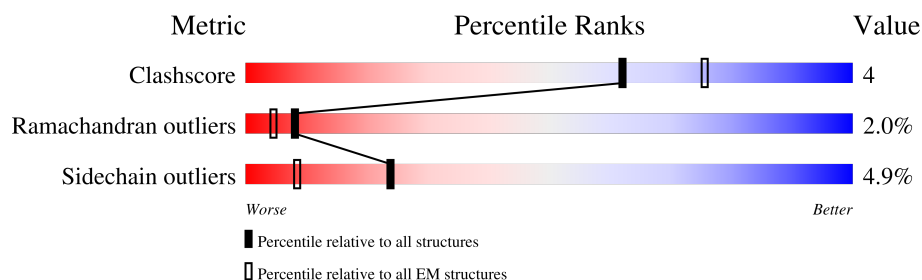
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 11.00 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.



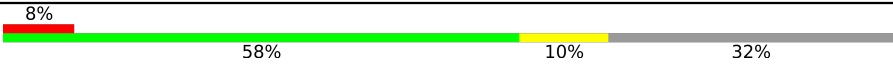
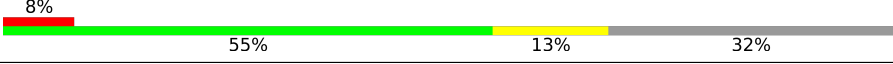
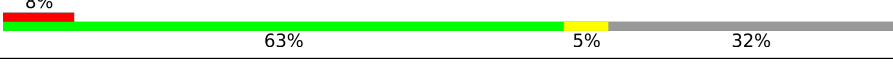
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	97 (10.50 - 11.50)

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	4387	9% 70% 11% • 19%
1	C	4387	9% 69% 11% • 19%
1	E	4387	9% 69% 11% • 19%
1	G	4387	9% 69% 11% • 19%

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Mol	Chain	Length	Quality of chain
2	B	158	
2	D	158	
2	F	158	
2	H	158	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 112936 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 Ryanodine Receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	3570	Total	C	N	O	S	0	0
			27416	17427	4732	5102	155		
1	C	3570	Total	C	N	O	S	0	0
			27416	17427	4732	5102	155		
1	E	3570	Total	C	N	O	S	0	0
			27416	17427	4732	5102	155		
1	G	3570	Total	C	N	O	S	0	0
			27416	17427	4732	5102	155		

- Molecule 2 is a protein called Peptidyl-prolyl cis-trans isomerase FKBP1B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	107	Total	C	N	O	S	0	0
			818	516	144	154	4		
2	D	107	Total	C	N	O	S	0	0
			818	516	144	154	4		
2	F	107	Total	C	N	O	S	0	0
			818	516	144	154	4		
2	H	107	Total	C	N	O	S	0	0
			818	516	144	154	4		

There are 204 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-50	MET	-	expression tag	UNP P68106
B	-49	ASN	-	expression tag	UNP P68106
B	-48	HIS	-	expression tag	UNP P68106
B	-47	LYS	-	expression tag	UNP P68106
B	-46	VAL	-	expression tag	UNP P68106
B	-45	HIS	-	expression tag	UNP P68106
B	-44	HIS	-	expression tag	UNP P68106
B	-43	HIS	-	expression tag	UNP P68106
B	-42	HIS	-	expression tag	UNP P68106

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-41	HIS	-	expression tag	UNP P68106
B	-40	HIS	-	expression tag	UNP P68106
B	-39	MET	-	expression tag	UNP P68106
B	-38	ASP	-	expression tag	UNP P68106
B	-37	GLU	-	expression tag	UNP P68106
B	-36	LYS	-	expression tag	UNP P68106
B	-35	THR	-	expression tag	UNP P68106
B	-34	THR	-	expression tag	UNP P68106
B	-33	GLY	-	expression tag	UNP P68106
B	-32	TRP	-	expression tag	UNP P68106
B	-31	ARG	-	expression tag	UNP P68106
B	-30	GLY	-	expression tag	UNP P68106
B	-29	GLY	-	expression tag	UNP P68106
B	-28	HIS	-	expression tag	UNP P68106
B	-27	VAL	-	expression tag	UNP P68106
B	-26	VAL	-	expression tag	UNP P68106
B	-25	GLU	-	expression tag	UNP P68106
B	-24	GLY	-	expression tag	UNP P68106
B	-23	LEU	-	expression tag	UNP P68106
B	-22	ALA	-	expression tag	UNP P68106
B	-21	GLY	-	expression tag	UNP P68106
B	-20	GLU	-	expression tag	UNP P68106
B	-19	LEU	-	expression tag	UNP P68106
B	-18	GLU	-	expression tag	UNP P68106
B	-17	GLN	-	expression tag	UNP P68106
B	-16	LEU	-	expression tag	UNP P68106
B	-15	ARG	-	expression tag	UNP P68106
B	-14	ALA	-	expression tag	UNP P68106
B	-13	ARG	-	expression tag	UNP P68106
B	-12	LEU	-	expression tag	UNP P68106
B	-11	GLU	-	expression tag	UNP P68106
B	-10	HIS	-	expression tag	UNP P68106
B	-9	HIS	-	expression tag	UNP P68106
B	-8	PRO	-	expression tag	UNP P68106
B	-7	GLN	-	expression tag	UNP P68106
B	-6	GLY	-	expression tag	UNP P68106
B	-5	GLN	-	expression tag	UNP P68106
B	-4	ARG	-	expression tag	UNP P68106
B	-3	GLU	-	expression tag	UNP P68106
B	-2	PRO	-	expression tag	UNP P68106
B	-1	GLU	-	expression tag	UNP P68106
B	0	LEU	-	expression tag	UNP P68106

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-50	MET	-	expression tag	UNP P68106
D	-49	ASN	-	expression tag	UNP P68106
D	-48	HIS	-	expression tag	UNP P68106
D	-47	LYS	-	expression tag	UNP P68106
D	-46	VAL	-	expression tag	UNP P68106
D	-45	HIS	-	expression tag	UNP P68106
D	-44	HIS	-	expression tag	UNP P68106
D	-43	HIS	-	expression tag	UNP P68106
D	-42	HIS	-	expression tag	UNP P68106
D	-41	HIS	-	expression tag	UNP P68106
D	-40	HIS	-	expression tag	UNP P68106
D	-39	MET	-	expression tag	UNP P68106
D	-38	ASP	-	expression tag	UNP P68106
D	-37	GLU	-	expression tag	UNP P68106
D	-36	LYS	-	expression tag	UNP P68106
D	-35	THR	-	expression tag	UNP P68106
D	-34	THR	-	expression tag	UNP P68106
D	-33	GLY	-	expression tag	UNP P68106
D	-32	TRP	-	expression tag	UNP P68106
D	-31	ARG	-	expression tag	UNP P68106
D	-30	GLY	-	expression tag	UNP P68106
D	-29	GLY	-	expression tag	UNP P68106
D	-28	HIS	-	expression tag	UNP P68106
D	-27	VAL	-	expression tag	UNP P68106
D	-26	VAL	-	expression tag	UNP P68106
D	-25	GLU	-	expression tag	UNP P68106
D	-24	GLY	-	expression tag	UNP P68106
D	-23	LEU	-	expression tag	UNP P68106
D	-22	ALA	-	expression tag	UNP P68106
D	-21	GLY	-	expression tag	UNP P68106
D	-20	GLU	-	expression tag	UNP P68106
D	-19	LEU	-	expression tag	UNP P68106
D	-18	GLU	-	expression tag	UNP P68106
D	-17	GLN	-	expression tag	UNP P68106
D	-16	LEU	-	expression tag	UNP P68106
D	-15	ARG	-	expression tag	UNP P68106
D	-14	ALA	-	expression tag	UNP P68106
D	-13	ARG	-	expression tag	UNP P68106
D	-12	LEU	-	expression tag	UNP P68106
D	-11	GLU	-	expression tag	UNP P68106
D	-10	HIS	-	expression tag	UNP P68106
D	-9	HIS	-	expression tag	UNP P68106

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-8	PRO	-	expression tag	UNP P68106
D	-7	GLN	-	expression tag	UNP P68106
D	-6	GLY	-	expression tag	UNP P68106
D	-5	GLN	-	expression tag	UNP P68106
D	-4	ARG	-	expression tag	UNP P68106
D	-3	GLU	-	expression tag	UNP P68106
D	-2	PRO	-	expression tag	UNP P68106
D	-1	GLU	-	expression tag	UNP P68106
D	0	LEU	-	expression tag	UNP P68106
F	-50	MET	-	expression tag	UNP P68106
F	-49	ASN	-	expression tag	UNP P68106
F	-48	HIS	-	expression tag	UNP P68106
F	-47	LYS	-	expression tag	UNP P68106
F	-46	VAL	-	expression tag	UNP P68106
F	-45	HIS	-	expression tag	UNP P68106
F	-44	HIS	-	expression tag	UNP P68106
F	-43	HIS	-	expression tag	UNP P68106
F	-42	HIS	-	expression tag	UNP P68106
F	-41	HIS	-	expression tag	UNP P68106
F	-40	HIS	-	expression tag	UNP P68106
F	-39	MET	-	expression tag	UNP P68106
F	-38	ASP	-	expression tag	UNP P68106
F	-37	GLU	-	expression tag	UNP P68106
F	-36	LYS	-	expression tag	UNP P68106
F	-35	THR	-	expression tag	UNP P68106
F	-34	THR	-	expression tag	UNP P68106
F	-33	GLY	-	expression tag	UNP P68106
F	-32	TRP	-	expression tag	UNP P68106
F	-31	ARG	-	expression tag	UNP P68106
F	-30	GLY	-	expression tag	UNP P68106
F	-29	GLY	-	expression tag	UNP P68106
F	-28	HIS	-	expression tag	UNP P68106
F	-27	VAL	-	expression tag	UNP P68106
F	-26	VAL	-	expression tag	UNP P68106
F	-25	GLU	-	expression tag	UNP P68106
F	-24	GLY	-	expression tag	UNP P68106
F	-23	LEU	-	expression tag	UNP P68106
F	-22	ALA	-	expression tag	UNP P68106
F	-21	GLY	-	expression tag	UNP P68106
F	-20	GLU	-	expression tag	UNP P68106
F	-19	LEU	-	expression tag	UNP P68106
F	-18	GLU	-	expression tag	UNP P68106

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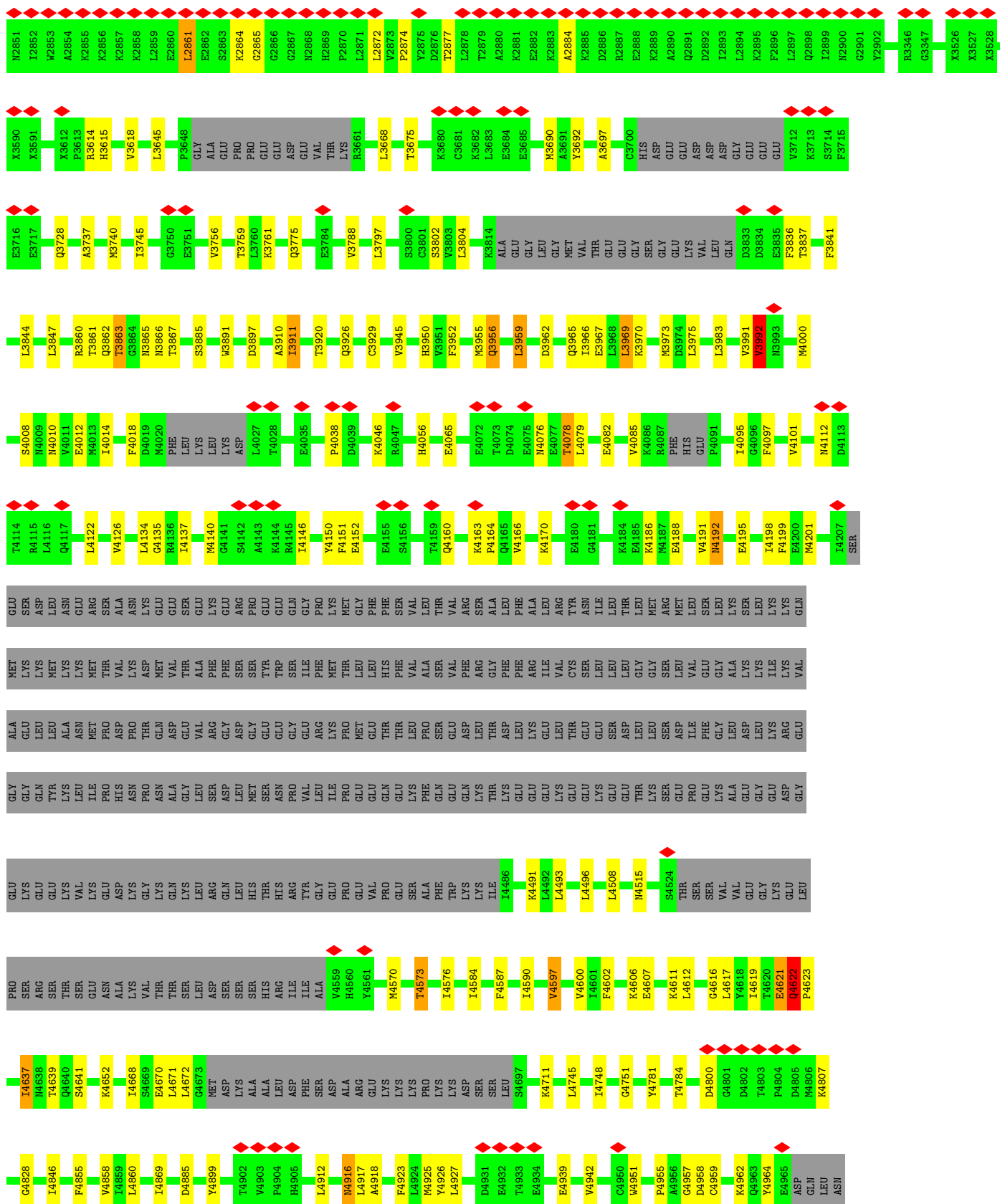
Chain	Residue	Modelled	Actual	Comment	Reference
F	-17	GLN	-	expression tag	UNP P68106
F	-16	LEU	-	expression tag	UNP P68106
F	-15	ARG	-	expression tag	UNP P68106
F	-14	ALA	-	expression tag	UNP P68106
F	-13	ARG	-	expression tag	UNP P68106
F	-12	LEU	-	expression tag	UNP P68106
F	-11	GLU	-	expression tag	UNP P68106
F	-10	HIS	-	expression tag	UNP P68106
F	-9	HIS	-	expression tag	UNP P68106
F	-8	PRO	-	expression tag	UNP P68106
F	-7	GLN	-	expression tag	UNP P68106
F	-6	GLY	-	expression tag	UNP P68106
F	-5	GLN	-	expression tag	UNP P68106
F	-4	ARG	-	expression tag	UNP P68106
F	-3	GLU	-	expression tag	UNP P68106
F	-2	PRO	-	expression tag	UNP P68106
F	-1	GLU	-	expression tag	UNP P68106
F	0	LEU	-	expression tag	UNP P68106
H	-50	MET	-	expression tag	UNP P68106
H	-49	ASN	-	expression tag	UNP P68106
H	-48	HIS	-	expression tag	UNP P68106
H	-47	LYS	-	expression tag	UNP P68106
H	-46	VAL	-	expression tag	UNP P68106
H	-45	HIS	-	expression tag	UNP P68106
H	-44	HIS	-	expression tag	UNP P68106
H	-43	HIS	-	expression tag	UNP P68106
H	-42	HIS	-	expression tag	UNP P68106
H	-41	HIS	-	expression tag	UNP P68106
H	-40	HIS	-	expression tag	UNP P68106
H	-39	MET	-	expression tag	UNP P68106
H	-38	ASP	-	expression tag	UNP P68106
H	-37	GLU	-	expression tag	UNP P68106
H	-36	LYS	-	expression tag	UNP P68106
H	-35	THR	-	expression tag	UNP P68106
H	-34	THR	-	expression tag	UNP P68106
H	-33	GLY	-	expression tag	UNP P68106
H	-32	TRP	-	expression tag	UNP P68106
H	-31	ARG	-	expression tag	UNP P68106
H	-30	GLY	-	expression tag	UNP P68106
H	-29	GLY	-	expression tag	UNP P68106
H	-28	HIS	-	expression tag	UNP P68106
H	-27	VAL	-	expression tag	UNP P68106

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-26	VAL	-	expression tag	UNP P68106
H	-25	GLU	-	expression tag	UNP P68106
H	-24	GLY	-	expression tag	UNP P68106
H	-23	LEU	-	expression tag	UNP P68106
H	-22	ALA	-	expression tag	UNP P68106
H	-21	GLY	-	expression tag	UNP P68106
H	-20	GLU	-	expression tag	UNP P68106
H	-19	LEU	-	expression tag	UNP P68106
H	-18	GLU	-	expression tag	UNP P68106
H	-17	GLN	-	expression tag	UNP P68106
H	-16	LEU	-	expression tag	UNP P68106
H	-15	ARG	-	expression tag	UNP P68106
H	-14	ALA	-	expression tag	UNP P68106
H	-13	ARG	-	expression tag	UNP P68106
H	-12	LEU	-	expression tag	UNP P68106
H	-11	GLU	-	expression tag	UNP P68106
H	-10	HIS	-	expression tag	UNP P68106
H	-9	HIS	-	expression tag	UNP P68106
H	-8	PRO	-	expression tag	UNP P68106
H	-7	GLN	-	expression tag	UNP P68106
H	-6	GLY	-	expression tag	UNP P68106
H	-5	GLN	-	expression tag	UNP P68106
H	-4	ARG	-	expression tag	UNP P68106
H	-3	GLU	-	expression tag	UNP P68106
H	-2	PRO	-	expression tag	UNP P68106
H	-1	GLU	-	expression tag	UNP P68106
H	0	LEU	-	expression tag	UNP P68106

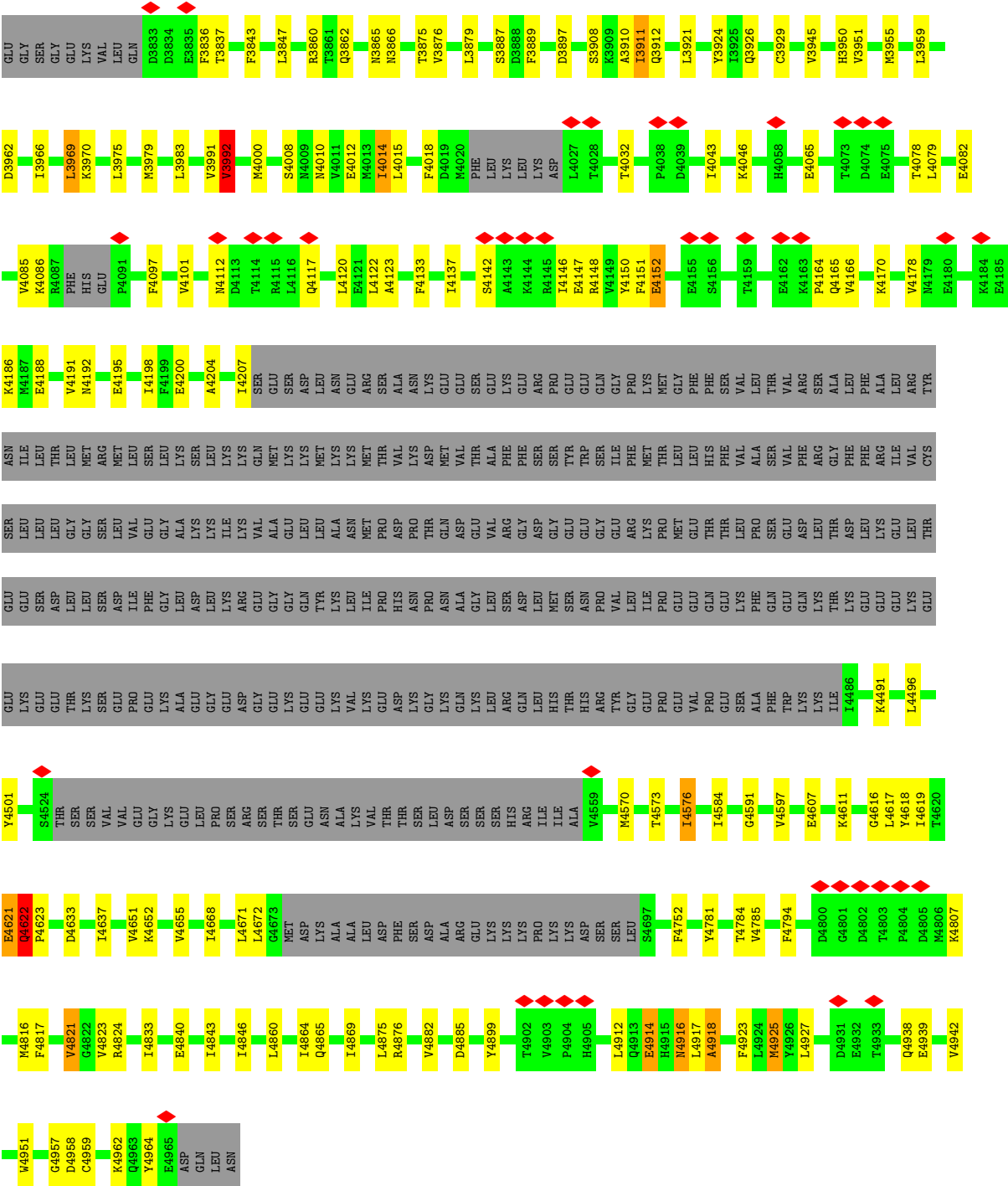




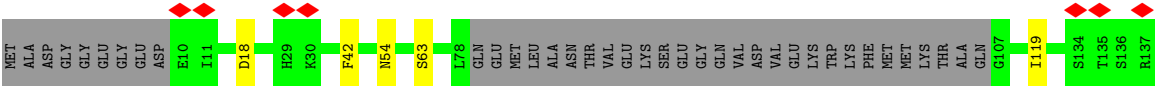
- Molecule 1: Ryanodine Receptor





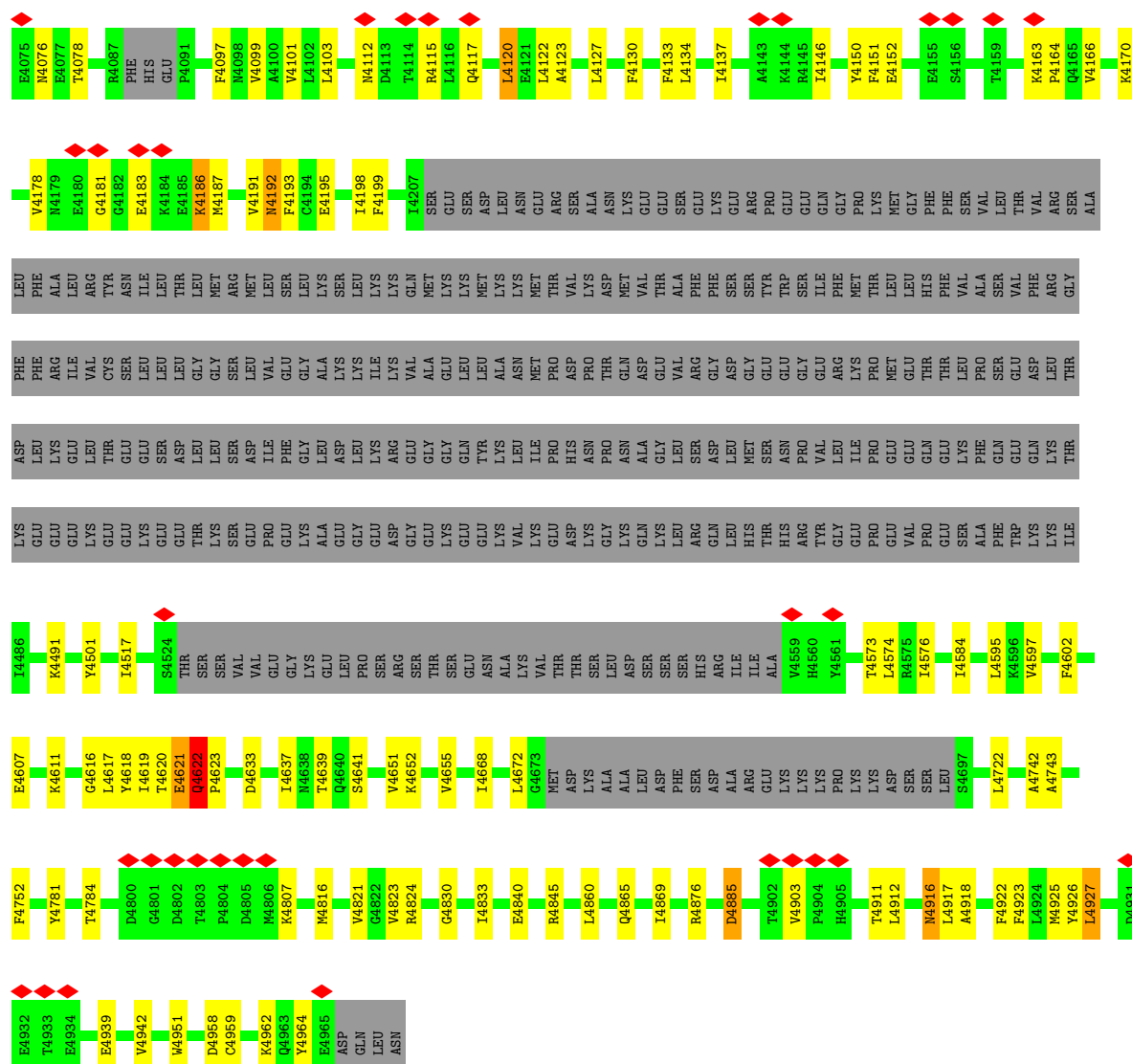


● Molecule 1: Ryanodine Receptor

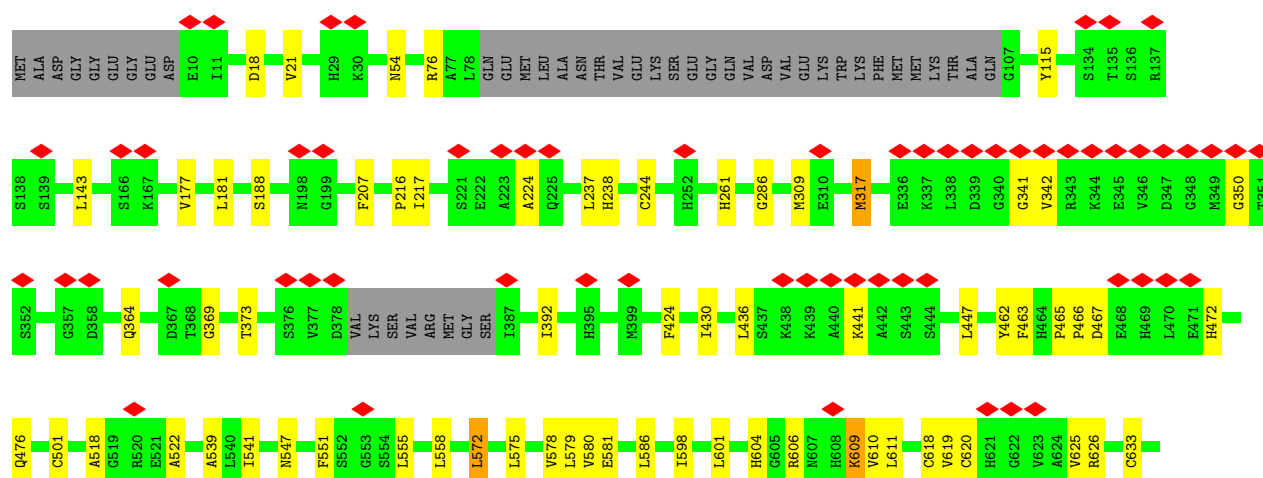








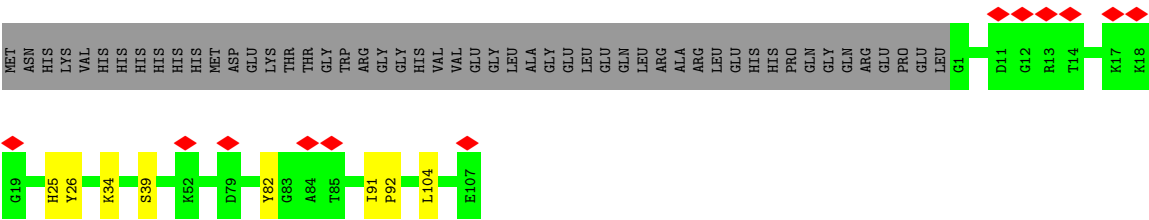
Molecule 1: Ryanodine Receptor







● Molecule 2: Peptidyl-prolyl cis-trans isomerase FKBP1B



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	13158	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	6000	Depositor
Magnification	62000	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.090	Depositor
Minimum map value	-0.046	Depositor
Average map value	-0.006	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.035	Depositor
Map size (Å)	536.0, 536.0, 536.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.34, 1.34, 1.34	Depositor

5 Model quality

5.1 Standard geometry

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	1.15	4/25831 (0.0%)	0.98	31/34865 (0.1%)
1	C	1.16	6/25831 (0.0%)	0.99	25/34865 (0.1%)
1	E	1.15	4/25831 (0.0%)	0.98	25/34865 (0.1%)
1	G	1.16	4/25831 (0.0%)	0.98	27/34865 (0.1%)
2	B	1.32	1/834 (0.1%)	1.02	2/1123 (0.2%)
2	D	1.33	0/834	1.00	0/1123
2	F	1.32	0/834	0.98	2/1123 (0.2%)
2	H	1.35	1/834 (0.1%)	1.01	0/1123
All	All	1.16	20/106660 (0.0%)	0.98	112/143952 (0.1%)

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	A	0	1
1	C	0	1
1	G	0	1
All	All	0	3

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	4183	GLU	CA-C	7.74	1.57	1.53
1	C	730	LEU	CA-C	6.51	1.55	1.52
1	G	4597	VAL	CA-CB	6.01	1.57	1.54
1	E	4597	VAL	CA-CB	5.85	1.57	1.54
1	C	1583	PRO	CA-C	5.81	1.57	1.52

The worst 5 of 112 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2248	MET	N-CA-C	7.79	119.85	111.36
1	A	1444	GLY	N-CA-C	7.72	121.92	110.42
1	C	2248	MET	N-CA-C	7.48	119.07	111.07
1	E	2248	MET	N-CA-C	7.26	119.27	111.36
1	G	2248	MET	N-CA-C	7.16	119.16	111.36

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	604	HIS	Peptide
1	C	3614	ARG	Peptide
1	G	3614	ARG	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	A	27416	0	25487	220	0
1	C	27416	0	25485	239	0
1	E	27416	0	25489	241	0
1	G	27416	0	25485	225	0
2	B	818	0	824	10	0
2	D	818	0	824	6	0
2	F	818	0	824	9	0
2	H	818	0	824	4	0
All	All	112936	0	105242	937	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

The worst 5 of 937 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:E:3945:VAL:HG11	1:E:3983:LEU:HD21	1.47	0.94
1:C:3778:MET:HE1	1:C:3847:LEU:HD22	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:874:LEU:HD23	1:A:878:LEU:HD13	1.50	0.94
1:G:4020:MET:HE3	1:G:4067:LEU:HD21	1.56	0.87
1:E:1172:THR:HG21	1:E:1190:LEU:HD13	1.58	0.86

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	3065/4387 (70%)	2784 (91%)	217 (7%)	64 (2%)	5	30
1	C	3065/4387 (70%)	2773 (90%)	236 (8%)	56 (2%)	6	34
1	E	3065/4387 (70%)	2765 (90%)	238 (8%)	62 (2%)	6	31
1	G	3065/4387 (70%)	2789 (91%)	211 (7%)	65 (2%)	5	30
2	B	105/158 (66%)	97 (92%)	7 (7%)	1 (1%)	12	49
2	D	105/158 (66%)	96 (91%)	8 (8%)	1 (1%)	12	49
2	F	105/158 (66%)	97 (92%)	6 (6%)	2 (2%)	6	32
2	H	105/158 (66%)	97 (92%)	8 (8%)	0	100	100
All	All	12680/18180 (70%)	11498 (91%)	931 (7%)	251 (2%)	8	31

5 of 251 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	472	HIS
1	A	1476	VAL
1	A	1549	SER
1	A	1779	SER
1	A	3897	ASP

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	2774/3480 (80%)	2654 (96%)	120 (4%)	26	47
1	C	2774/3480 (80%)	2629 (95%)	145 (5%)	21	42
1	E	2774/3480 (80%)	2628 (95%)	146 (5%)	20	41
1	G	2774/3480 (80%)	2631 (95%)	143 (5%)	21	42
2	B	88/131 (67%)	87 (99%)	1 (1%)	65	76
2	D	88/131 (67%)	87 (99%)	1 (1%)	65	76
2	F	88/131 (67%)	87 (99%)	1 (1%)	65	76
2	H	88/131 (67%)	87 (99%)	1 (1%)	65	76
All	All	11448/14444 (79%)	10890 (95%)	558 (5%)	24	43

5 of 558 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	1657	THR
1	G	1790	LYS
1	G	1644	LEU
1	G	3690	MET
1	C	2117	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 243 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	4581	HIS
1	G	2900	ASN
1	E	1002	ASN
1	G	2891	GLN
1	G	4638	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	27
1	G	27
1	E	27
1	A	27

The worst 5 of 108 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	3612:UNK	C	3613:PRO	N	76.52
1	G	3612:UNK	C	3613:PRO	N	75.29
1	E	3612:UNK	C	3613:PRO	N	74.52
1	A	3612:UNK	C	3613:PRO	N	74.10
1	E	2701:UNK	C	2702:PHE	N	43.29

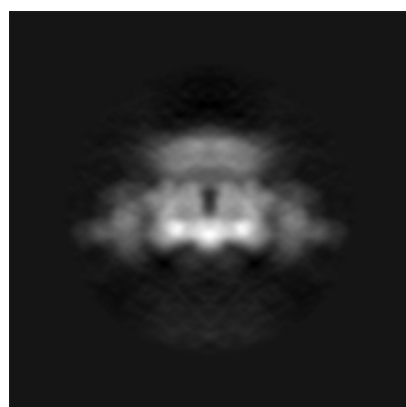
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8303. These allow visual inspection of the internal detail of the map and identification of artifacts.

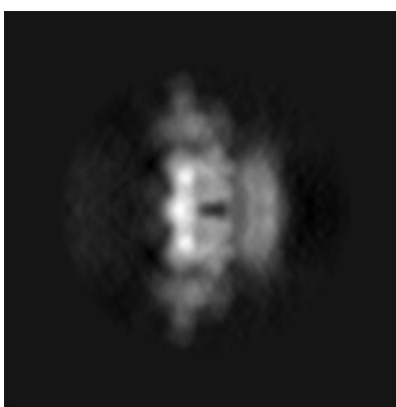
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

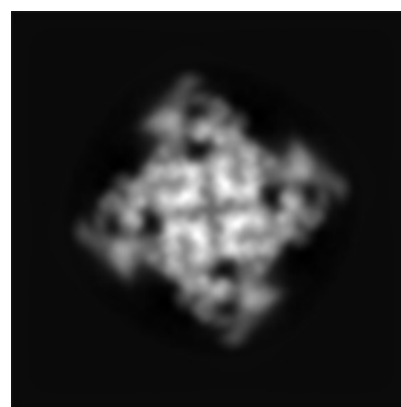
6.1.1 Primary map



X



Y

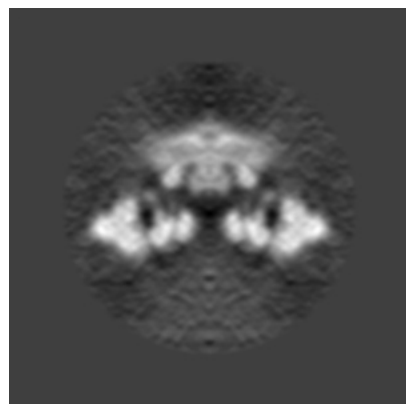


Z

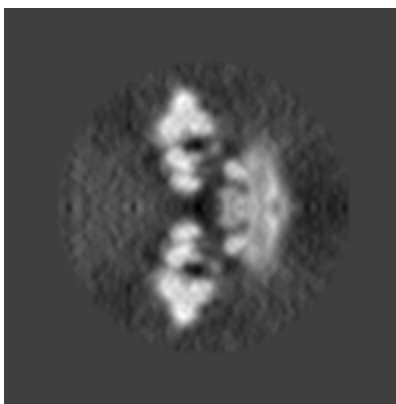
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

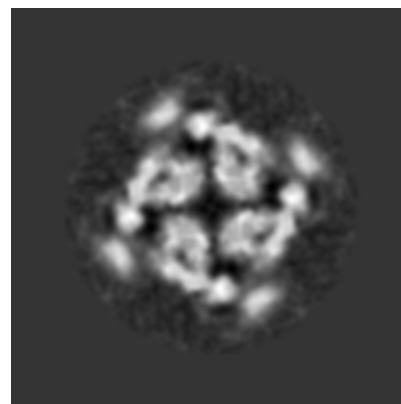
6.2.1 Primary map



X Index: 200



Y Index: 200

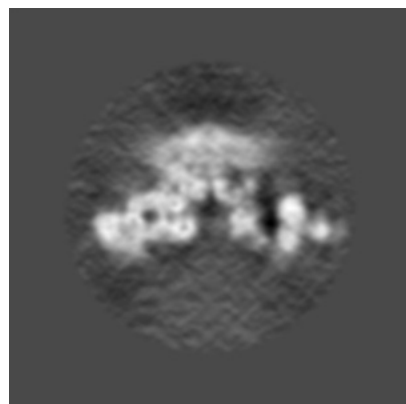


Z Index: 200

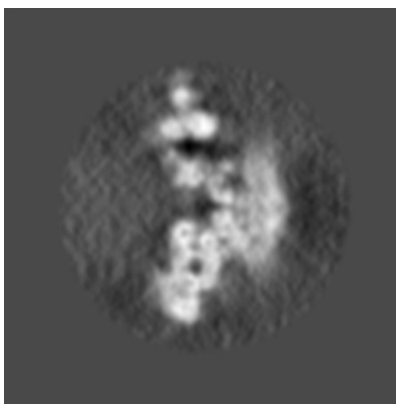
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

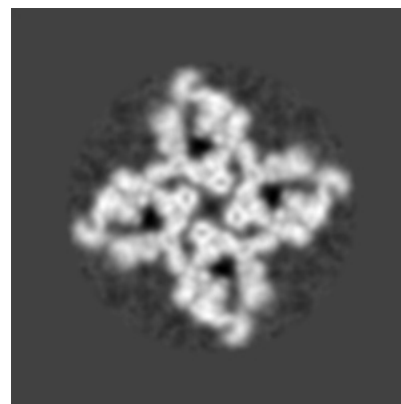
6.3.1 Primary map



X Index: 190



Y Index: 210

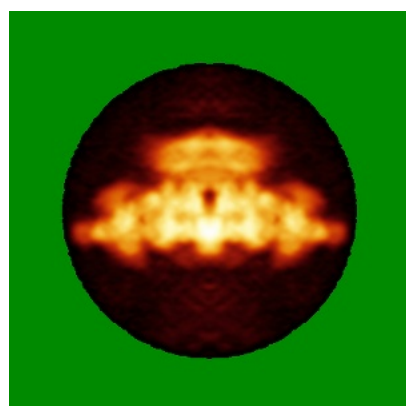


Z Index: 181

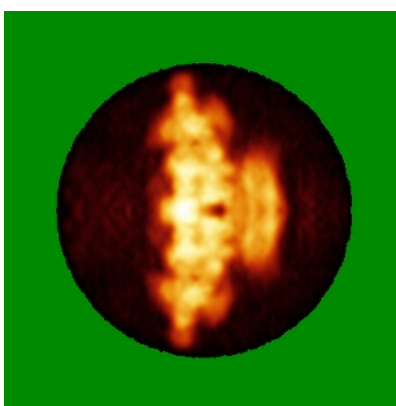
The images above show the largest variance slices of the map in three orthogonal directions.

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

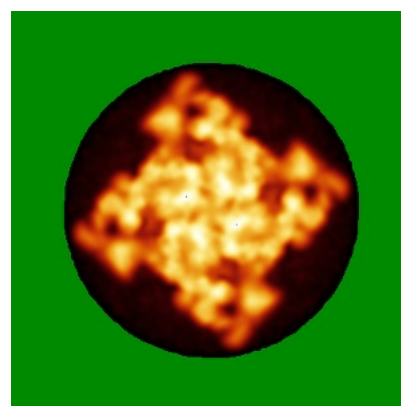
6.4.1 Primary map



X



Y

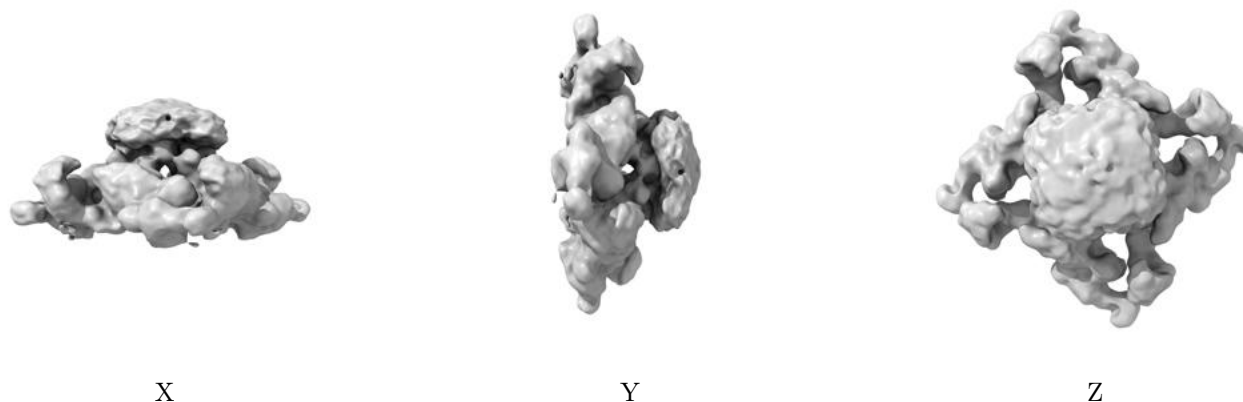


Z

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

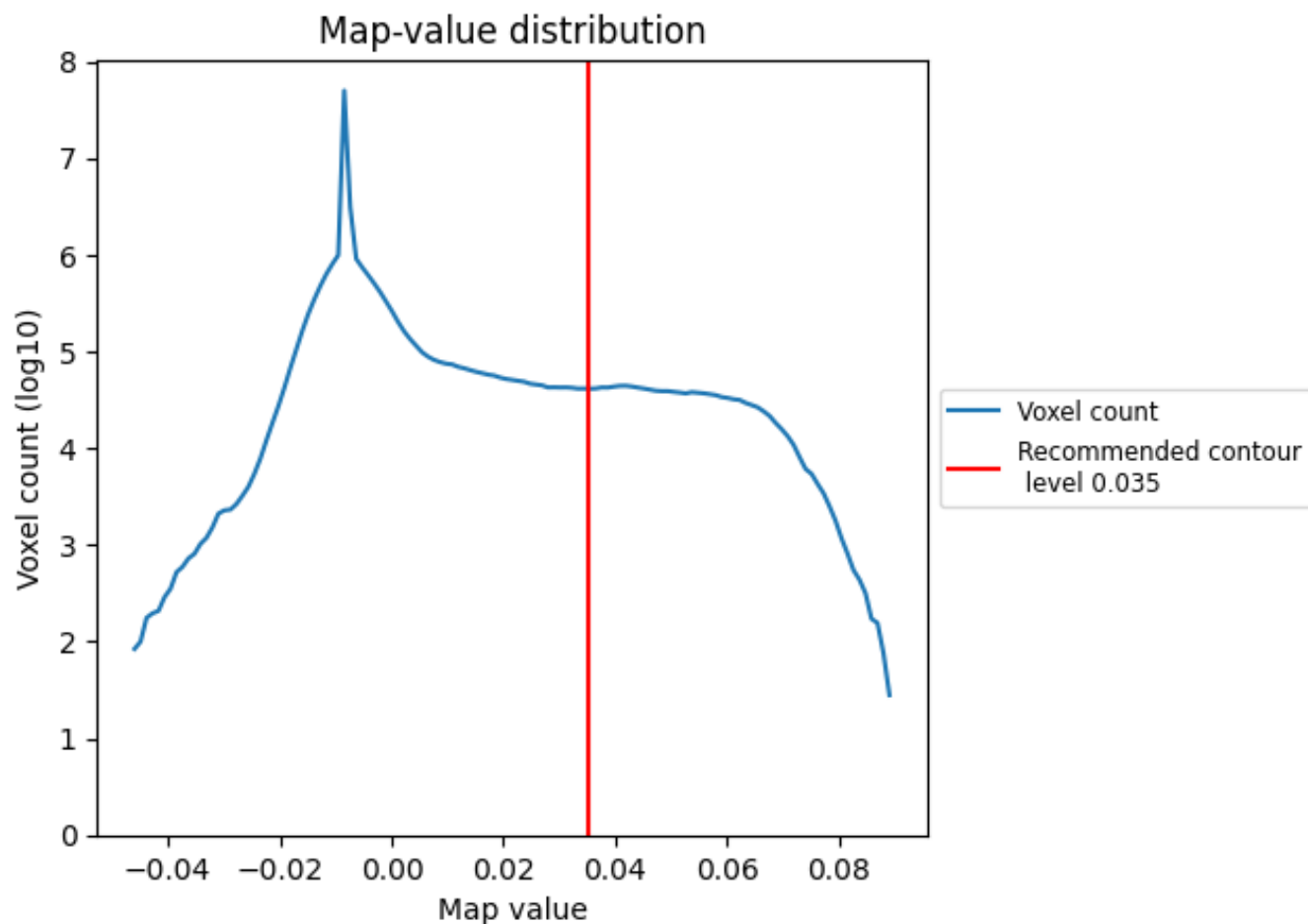
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

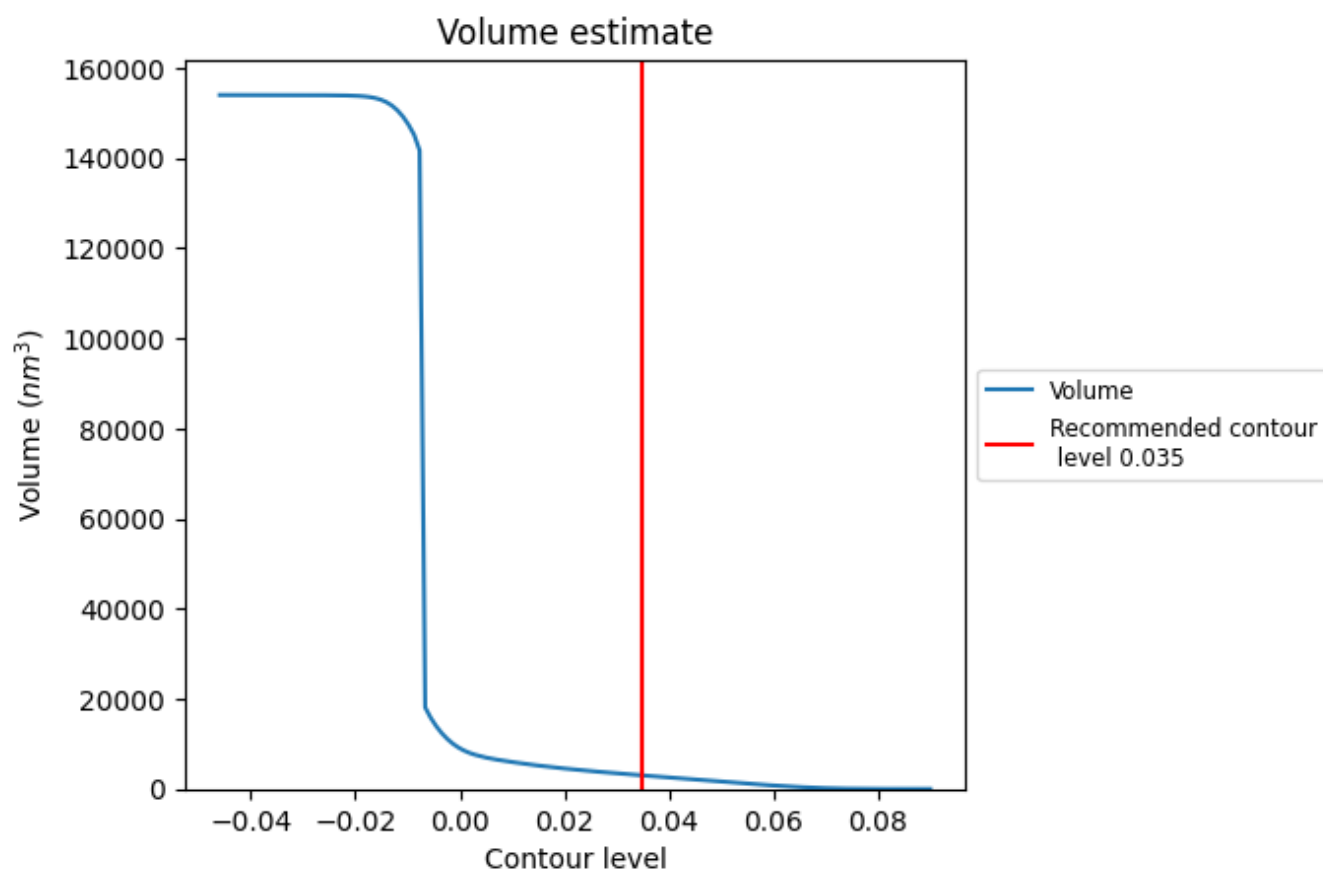
This section contains the results of statistical analysis of the map.

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.

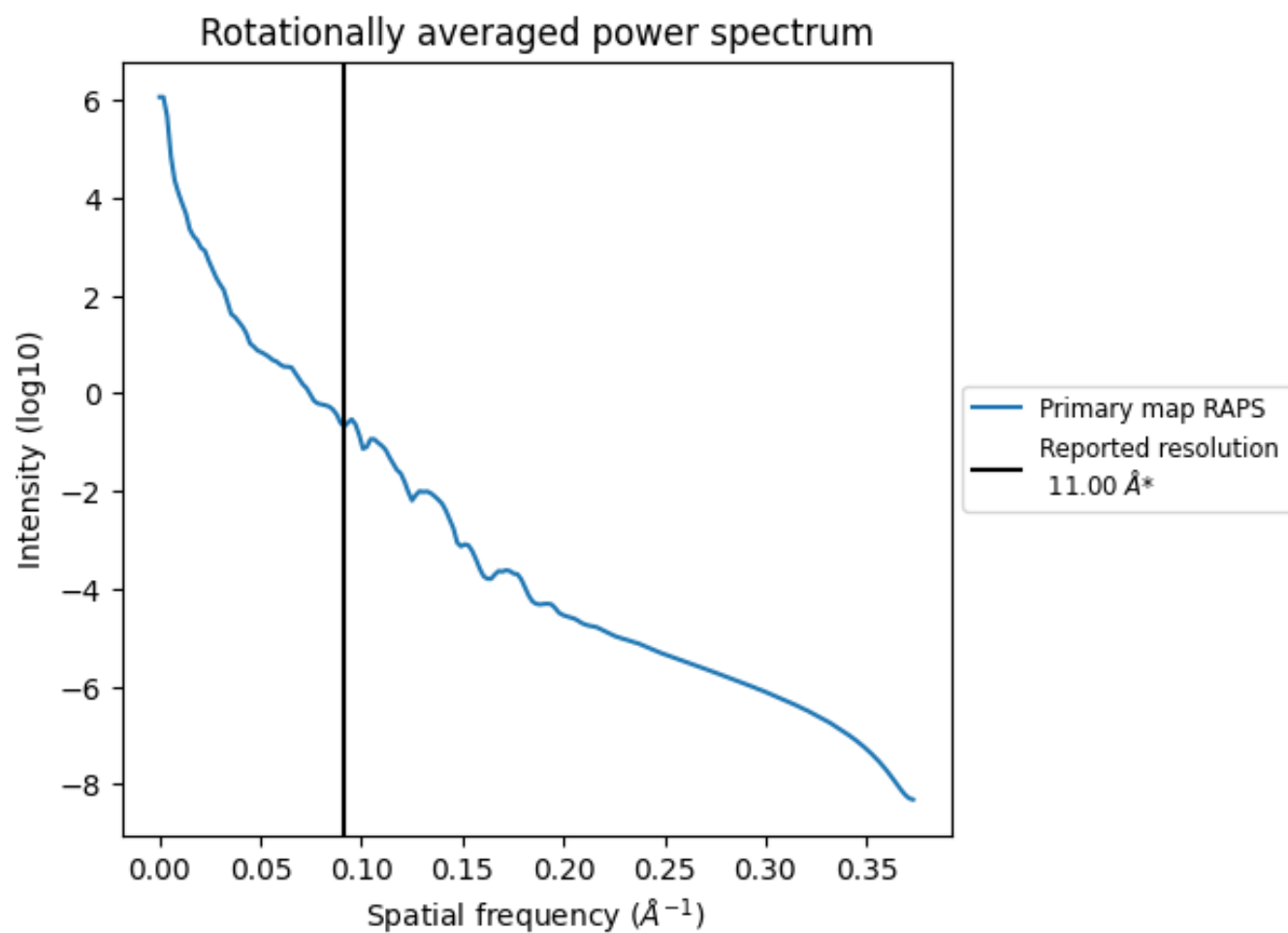
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 3008 nm³; this corresponds to an approximate mass of 2718 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 0.091 Å⁻¹

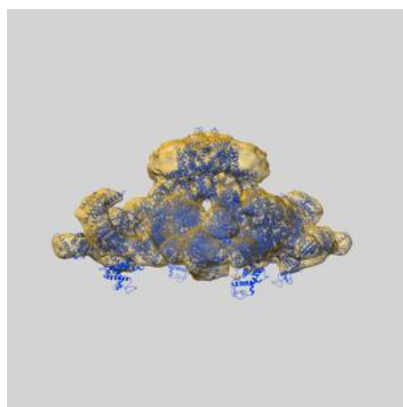
8 Fourier-Shell correlation

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

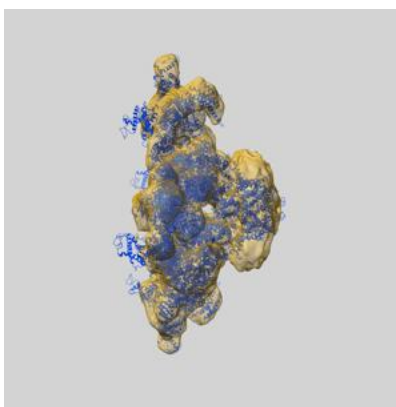
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8303 and PDB model 5L1D. Per-residue inclusion information can be found in section [3](#) on page [10](#).

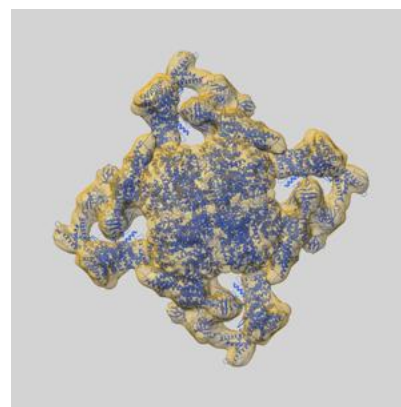
9.1 Map-model overlay [i](#)



X



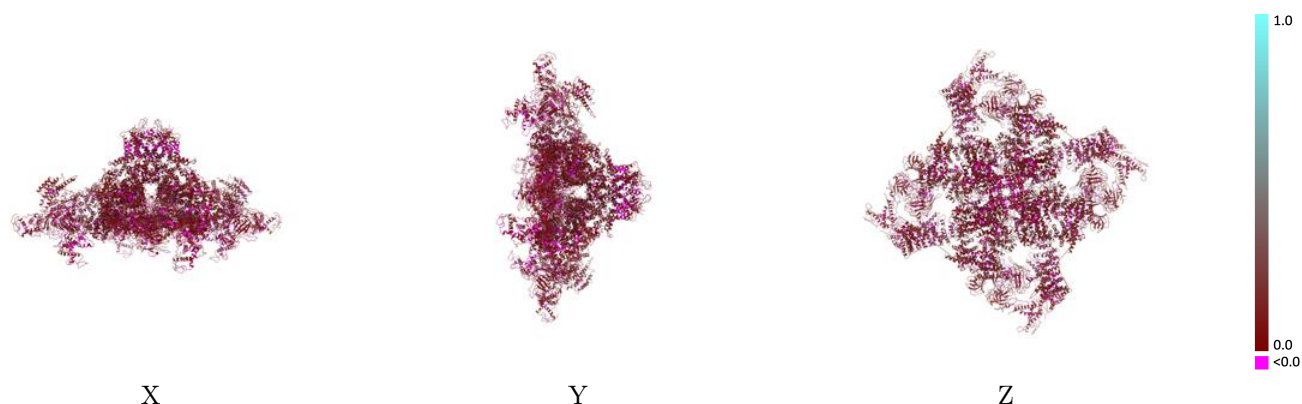
Y



Z

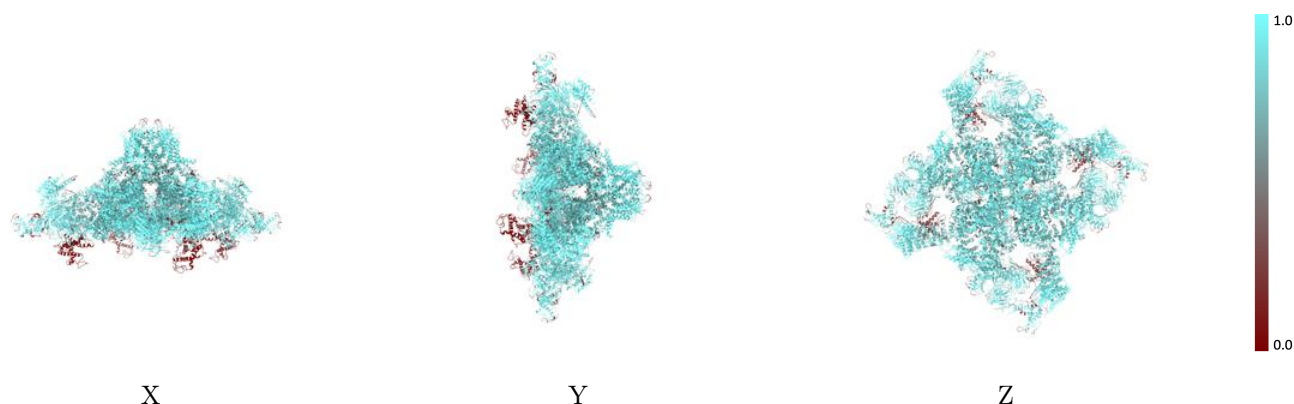
The images above show the 3D surface view of the map at the recommended contour level 0.035 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



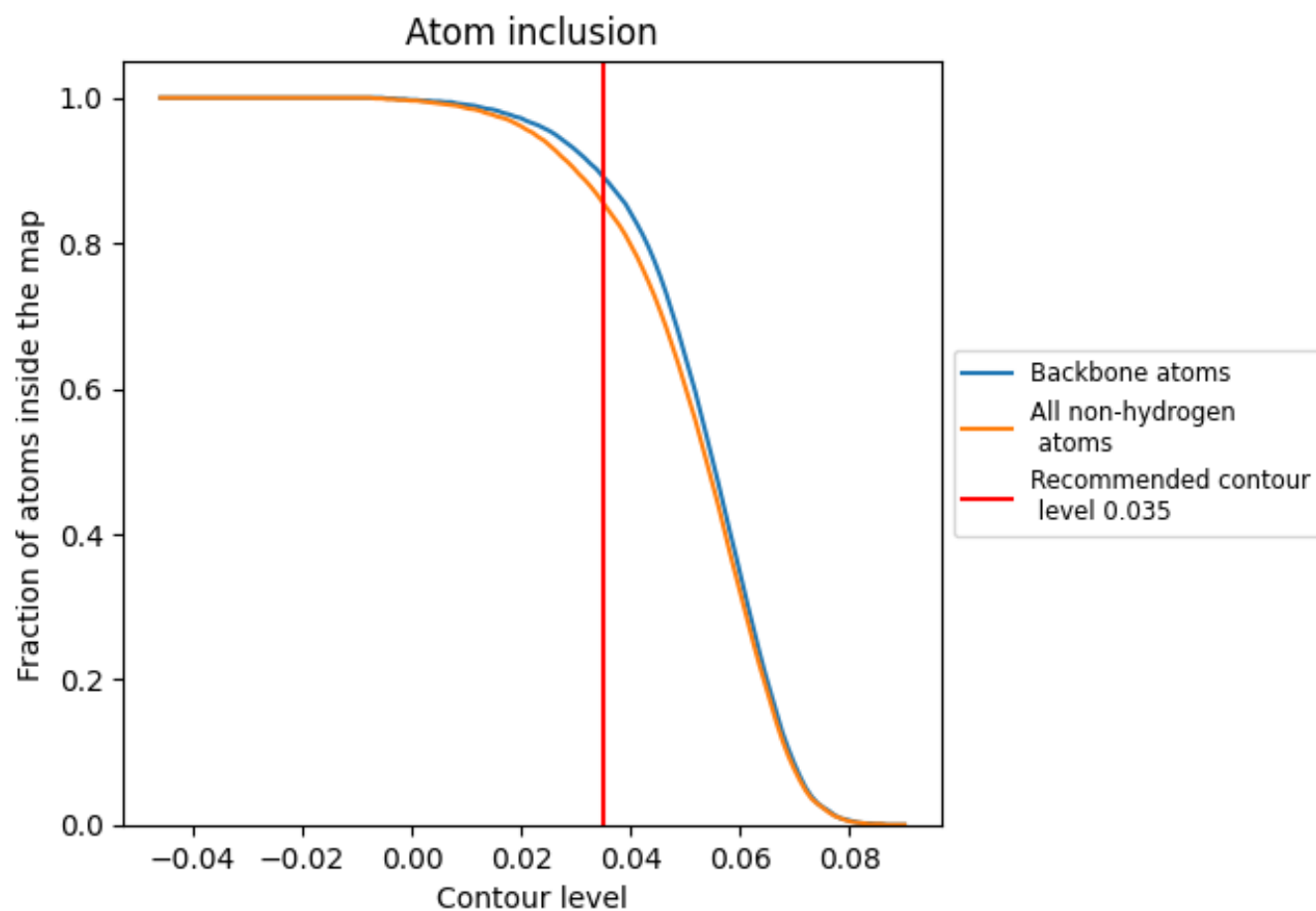
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.035).

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 86% 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 (0.035) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8550	0.0820
A	0.8560	0.0820
B	0.8470	0.0660
C	0.8560	0.0820
D	0.8460	0.0720
E	0.8550	0.0820
F	0.8440	0.0750
G	0.8560	0.0830
H	0.8380	0.0700

1.0

0.0

<0.0