



wwPDB EM Validation Summary Report ⓘ

Mar 22, 2026 – 09:13 PM UTC

PDB ID : 4BOI / pdb_00004boi
EMDB ID : EMD-2377
Title : The structure and super-organization of acetylcholine receptor-rapsyn complexes class A
Authors : Zuber, B.; Unwin, N.
Deposited on : 2013-05-20
Resolution : 41.00 Å(reported)
Based on initial model : 2BG9

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 : **NOT EXECUTED**
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

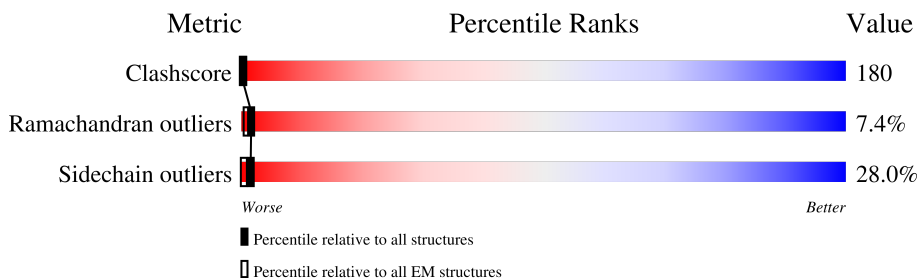
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 41.00 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	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

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	461	18% 6% 40% 29% 6% 20%
1	D	461	12% 5% 44% 25% 5% 20%
2	B	493	8% • 42% 23% 5% 25%
3	C	522	10% 6% 36% 23% 5% 29%
4	E	505	7% 5% 38% 25% 5% 27%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14924 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 ACETYLCHOLINE RECEPTOR SUBUNIT ALPHA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	370	Total	C	N	O	S	0	0
			2991	1954	478	540	19		
1	D	370	Total	C	N	O	S	0	0
			2991	1954	478	540	19		

- Molecule 2 is a protein called ACETYLCHOLINE RECEPTOR BETA SUBUNIT.

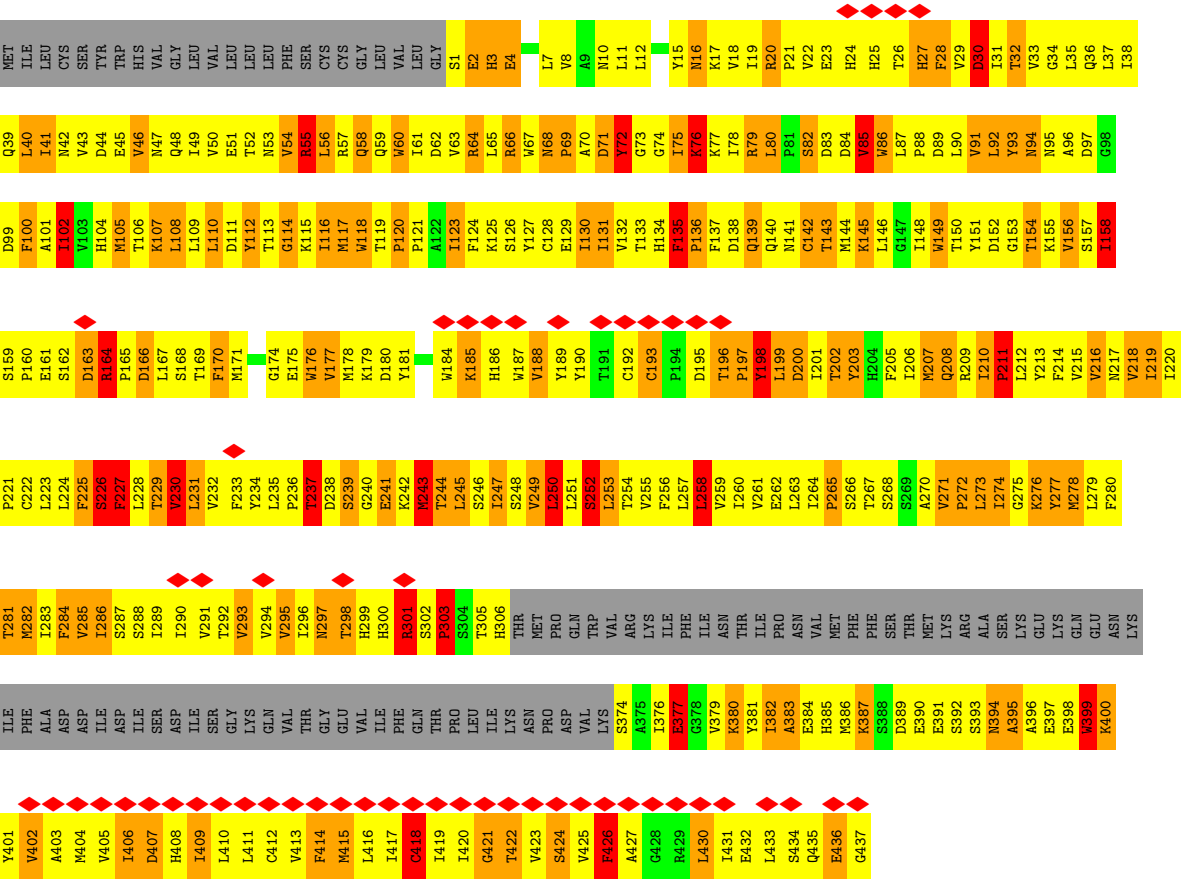
Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	370	Total	C	N	O	S	0	0
			2972	1938	465	554	15		

- Molecule 3 is a protein called ACETYLCHOLINE RECEPTOR DELTA SUBUNIT.

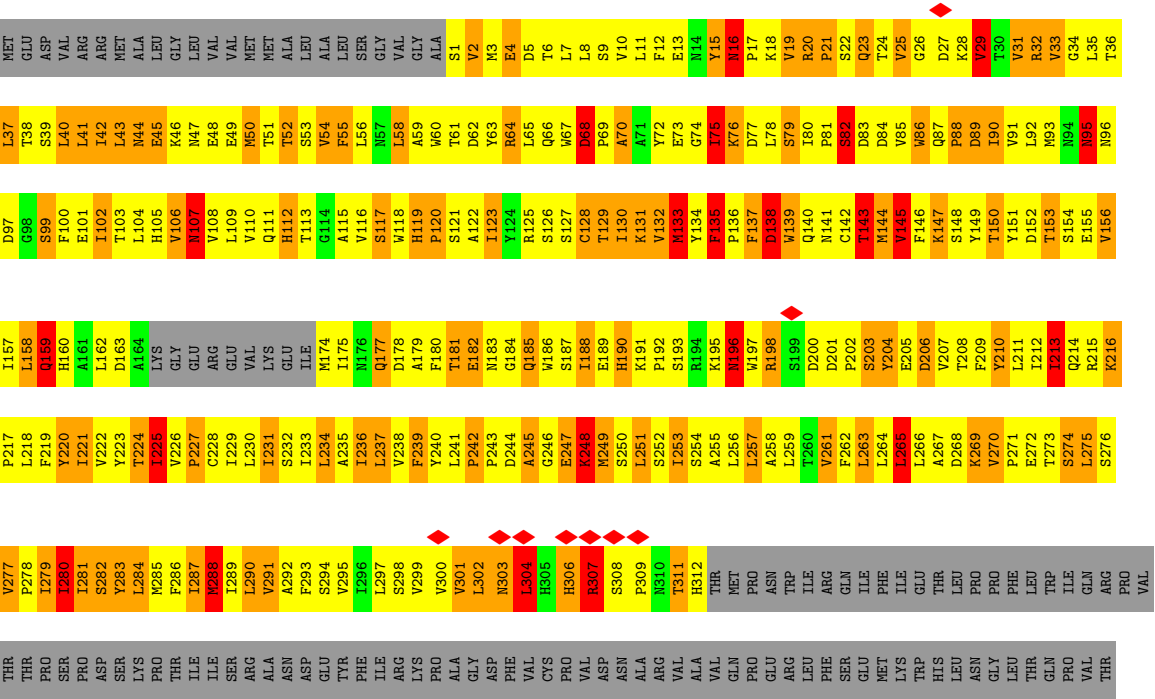
Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	370	Total	C	N	O	S	0	1
			2983	1944	489	536	14		

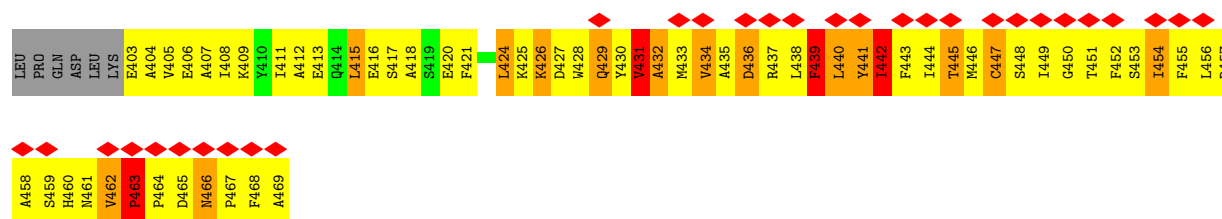
- Molecule 4 is a protein called ACETYLCHOLINE RECEPTOR GAMMA SUBUNIT.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	371	Total	C	N	O	S	0	1
			2987	1948	478	551	10		

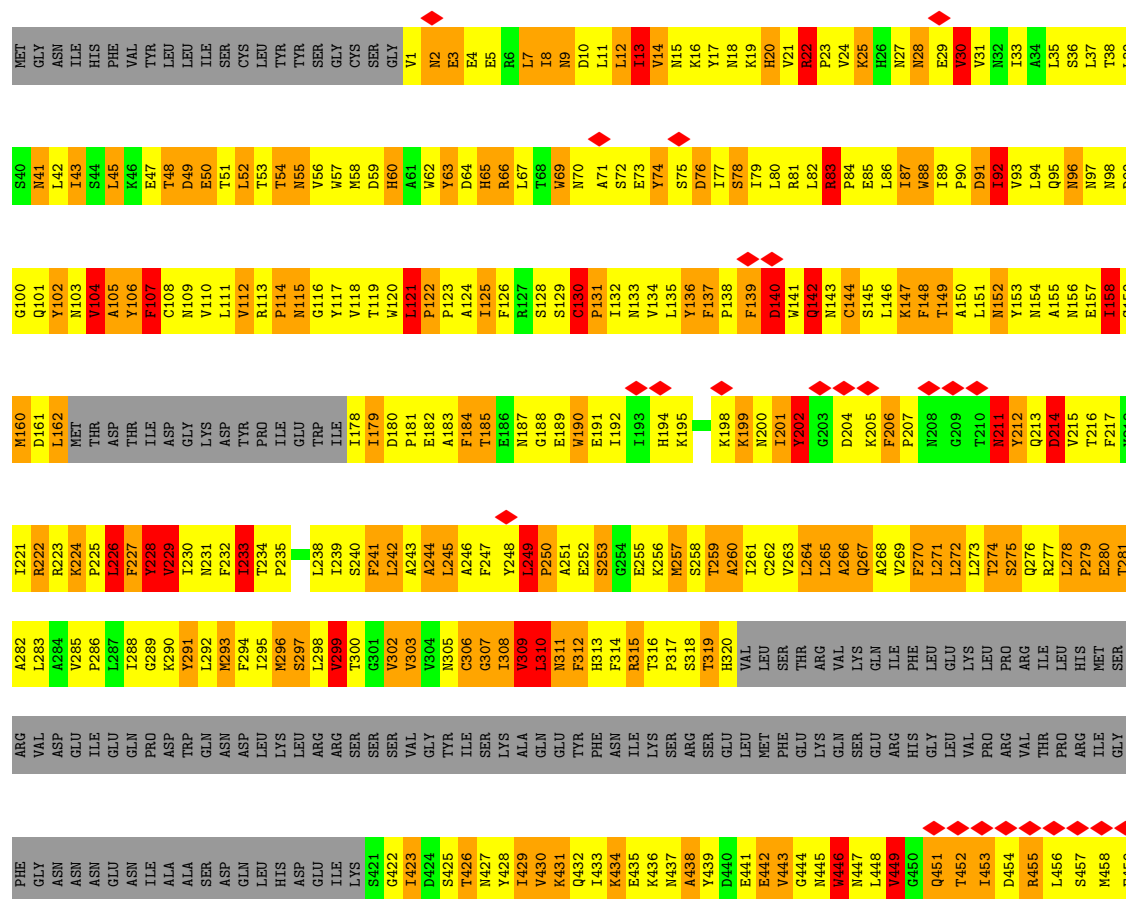


● Molecule 2: ACETYLCHOLINE RECEPTOR BETA SUBUNIT

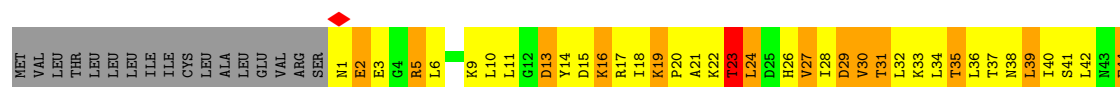




• Molecule 3: ACETYLCHOLINE RECEPTOR DELTA SUBUNIT



• Molecule 4: ACETYLCHOLINE RECEPTOR GAMMA SUBUNIT



T465	F466	L467	T468	G469	H470	L471	M472	Q473	V474	P475	E476	F477	PRO	PHE	PRO	GLY	ASP	PRO	ARG	LYS	TYR	VAL	PRO																																				
LEU	ALA	ASN	PHE	ALA	PRO	GLU	ILE	SER	SER	PHE	GLY	ILE	ILE	LYS	ALA	GLU	TYR	ILE	LEU	LYS	PRO	ARG	SER	GLU	GLN	LYS	ASP	ARG	VAL	ASN	LYS	MET	HIS	LEU	PHE	LEU	PRO	LYS	TYR	LEU	GLY	MET	HIS	LEU	GLU	PRO	GLU												
Y285	L286	F287	V288	M290	F291	V292	S293	L294	V295	I296	V297	T298	N299	C300	I301	V302	V303	L304	N305	V306	S307	L308	R309	T310	P311	N312	T313	H314	SER	LEU	SER	GLU	LYS	LYS	ILE	LYS	HIS	LEU	PHE	LEU	PHE	LEU	PRO	LYS	TYR	LEU	GLY	MET	HIS	LEU	GLU	PRO	GLU						
I225	I226	A227	P228	C229	V230	L231	I232	S233	S234	L235	V236	V237	L238	V239	F241	L242	P243	A244	Q245	A246	G247	C248	Q249	K250	C251	T252	L253	S254	I255	S256	V257	L258	L259	A260	Q261	T262	I263	F264	L265	F266	L267	I268	A269	Q270	F271	L272	P273	E274	T275	S276	L277	V278	V279	P280	L281	I282	G283	K284	
GLU	VAL	VAL	GLU	TRP	ILE	HIS	I172	D173	P174	E175	D176	F177	T178	E179	N180	A181	E182	W183	T184	I185	R186	H187	R188	P189	A190	K191	T192	N193	Y194	N195	W196	L198	T199	K200	D201	D202	I203	D204	F205	Q206	E207	I208	I209	F210	F211	L212	I213	I214	Q215	R216	K217	P218	L219	F220	Y221	I222	G223	N224	
A105	E106	V107	L108	Y109	T110	N111	D112	G113	S114	M115	Y116	W117	L118	P119	P120	A121	I122	Y123	R124	S125	T126	C127	P128	I129	A130	V131	T132	Y133	F134	P135	F136	D137	W138	Q139	N140	C141	S142	L143	V144	F145	R146	S147	Q148	T149	Y150	N151	A152	H153	E154	V155	N156	L157	Q158	F159	E160	A161	E162	E163	G164
K45	E46	E47	A48	L49	T50	T51	N52	V53	W54	I55	E56	I57	Q58	W59	N60	D61	Y62	R63	L64	S65	W66	N67	T68	S69	E70	Y71	E72	G73	I74	D75	L76	V77	R78	I79	P80	S81	E82	L83	L84	W85	L86	P87	D88	V89	V90	L91	E92	R93	N94	V95	D96	G97	Q98	F99	V100	V101	A102	Y103	Y104

4 Experimental information

Property	Value	Source
EM reconstruction method	TOMOGRAPHY	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of tilted images used	3564	Depositor
Resolution determination method	Not provided	
CTF correction method	Not provided	
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	3000	Depositor
Maximum defocus (nm)	6000	Depositor
Magnification	80213	Depositor
Image detector	GATAN ULTRASCAN 4000 (4k x 4k)	Depositor
Maximum voxel value	1.230	Depositor
Minimum voxel value	-0.688	Depositor
Average voxel value	0.000	Depositor
Voxel value standard deviation	0.078	Depositor
Recommended contour level	0.413	Depositor
Tomogram size ($\text{\AA}$)	448.8, 448.8, 448.8	wwPDB
Tomogram dimensions	60, 60, 60	wwPDB
Tomogram angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Grid spacing ($\text{\AA}$)	7.48, 7.48, 7.48	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	0.96	9/3069 (0.3%)	1.55	70/4186 (1.7%)
1	D	0.98	13/3069 (0.4%)	1.49	56/4186 (1.3%)
2	B	1.05	17/3048 (0.6%)	1.51	61/4162 (1.5%)
3	C	0.98	13/3059 (0.4%)	1.55	72/4175 (1.7%)
4	E	0.96	13/3057 (0.4%)	1.50	72/4174 (1.7%)
All	All	0.99	65/15302 (0.4%)	1.52	331/20883 (1.6%)

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	D	0	2
3	C	0	2
All	All	0	4

The worst 5 of 65 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	129	THR	C-N	-18.39	1.14	1.34
1	D	208	GLN	C-N	12.71	1.51	1.33
4	E	182	GLU	C-N	12.51	1.44	1.33
3	C	265	LEU	C-N	10.84	1.48	1.33
4	E	306	VAL	C-N	-8.82	1.22	1.33

The worst 5 of 331 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	122	PRO	CA-C-N	13.73	134.66	119.83
3	C	122	PRO	C-N-CA	13.73	134.66	119.83
3	C	460	ILE	N-CA-C	-13.48	100.98	111.90
4	E	238	LEU	N-CA-C	-13.46	96.08	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	441	TYR	N-CA-C	-12.79	97.03	112.89

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	63	TYR	Sidechain
3	C	74	TYR	Sidechain
1	D	277	TYR	Sidechain
1	D	72	TYR	Sidechain

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	2991	0	3006	1097	0
1	D	2991	0	3006	1091	0
2	B	2972	0	2954	1128	0
3	C	2983	0	2987	1203	0
4	E	2987	0	2994	1110	0
All	All	14924	0	14947	5364	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 180.

The worst 5 of 5364 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)
4:E:183:TRP:CB	4:E:216:ARG:HG2	1.33	1.51
2:B:134:TYR:CE1	2:B:213:ILE:HG13	1.44	1.49
1:A:167:LEU:HD12	1:A:178:MET:CB	1.43	1.45
1:A:167:LEU:CD1	1:A:178:MET:HB2	1.46	1.44
3:C:316:THR:CG2	3:C:317:PRO:HD2	1.53	1.38

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	A	366/461 (79%)	288 (79%)	50 (14%)	28 (8%)	1	10
1	D	366/461 (79%)	294 (80%)	41 (11%)	31 (8%)	0	9
2	B	364/493 (74%)	274 (75%)	58 (16%)	32 (9%)	0	8
3	C	364/522 (70%)	288 (79%)	58 (16%)	18 (5%)	1	16
4	E	365/505 (72%)	281 (77%)	58 (16%)	26 (7%)	1	11
All	All	1825/2442 (75%)	1425 (78%)	265 (14%)	135 (7%)	1	10

5 of 135 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	27	HIS
1	A	76	LYS
1	A	83	ASP
1	A	102	ILE

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	343/427 (80%)	236 (69%)	107 (31%)	0	2
1	D	343/427 (80%)	248 (72%)	95 (28%)	0	3
2	B	340/449 (76%)	261 (77%)	79 (23%)	1	5
3	C	335/475 (70%)	240 (72%)	95 (28%)	0	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	E	337/463 (73%)	238 (71%)	99 (29%)	0	2
All	All	1698/2241 (76%)	1223 (72%)	475 (28%)	1	3

5 of 475 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	206	PHE
4	E	253	LEU
1	D	28	PHE
4	E	232	ILE
4	E	473	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 71 such sidechains are listed below:

Mol	Chain	Res	Type
4	E	98	GLN
4	E	148	GLN
4	E	249	GLN
3	C	41	ASN
3	C	20	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
2	B	1

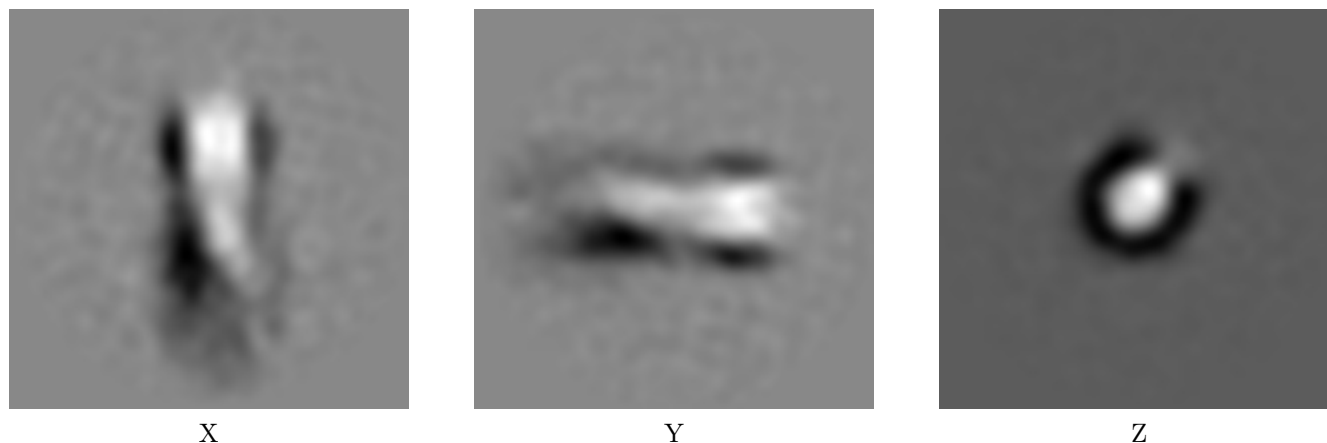
All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	129:THR	C	130:ILE	N	1.14

6 Tomogram visualisation [i](#)

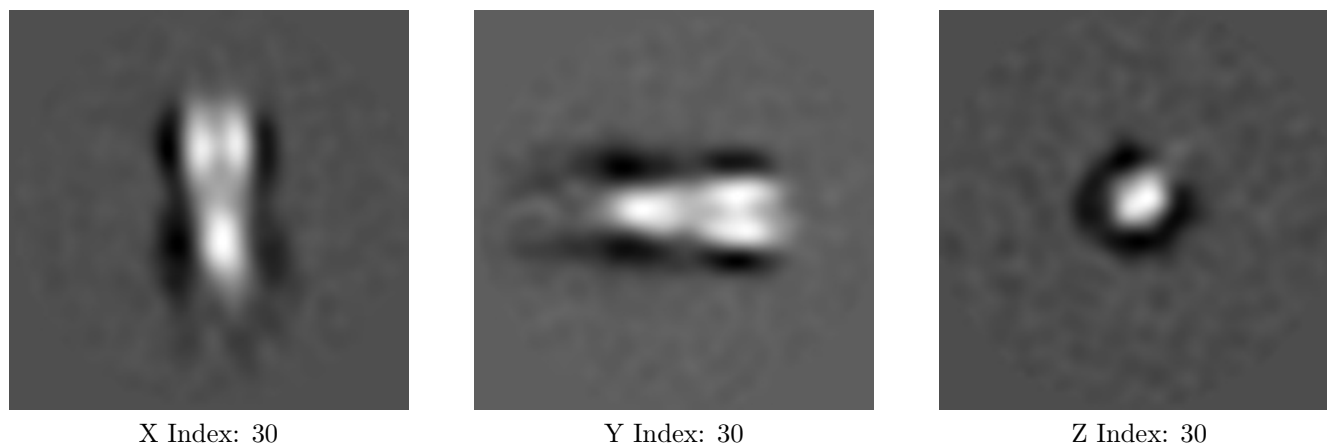
This section contains visualisations of the EMDB entry EMD-2377. These allow visual inspection of the internal detail of the tomogram and identification of artifacts.

6.1 Orthogonal projections [i](#)



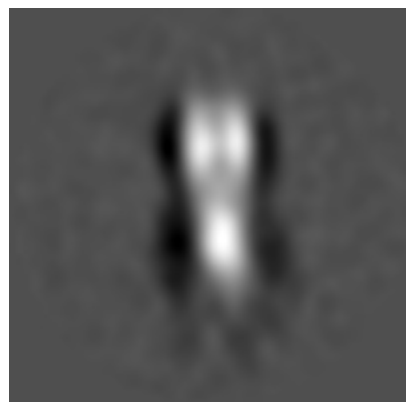
The images above show the tomogram projected in three orthogonal directions.

6.2 Central slices [i](#)

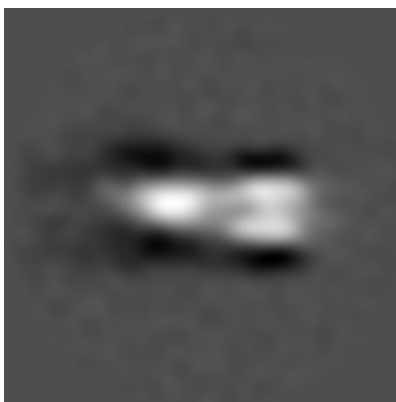


The images above show central slices of the tomogram in three orthogonal directions.

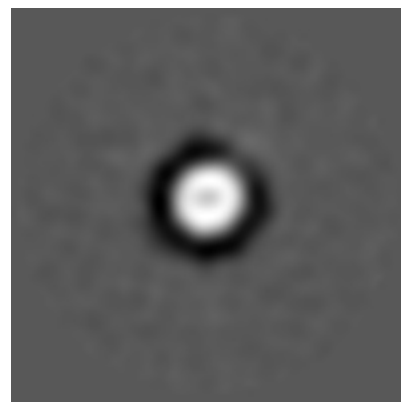
6.3 Largest variance slices [i](#)



X Index: 30



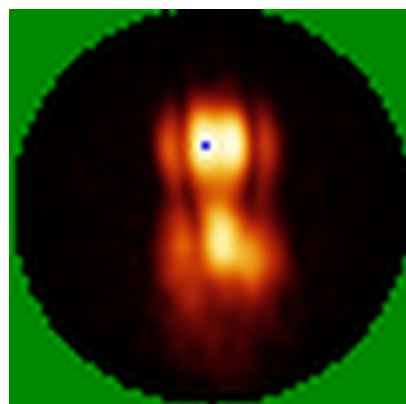
Y Index: 32



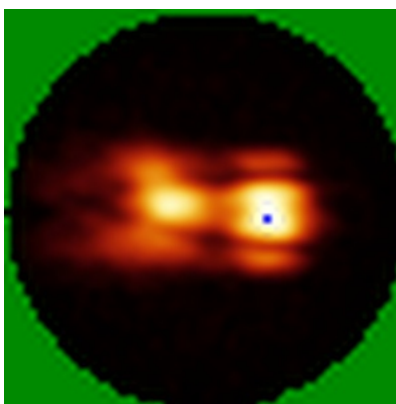
Z Index: 39

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

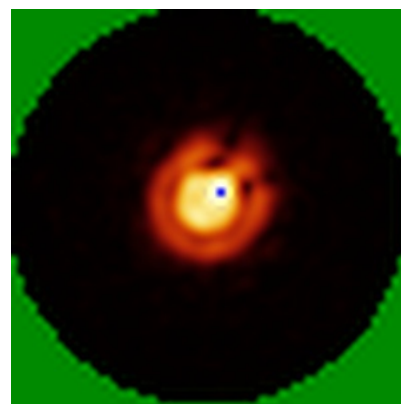
6.4 Orthogonal standard-deviation projections (False-color) [i](#)



X



Y



Z

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

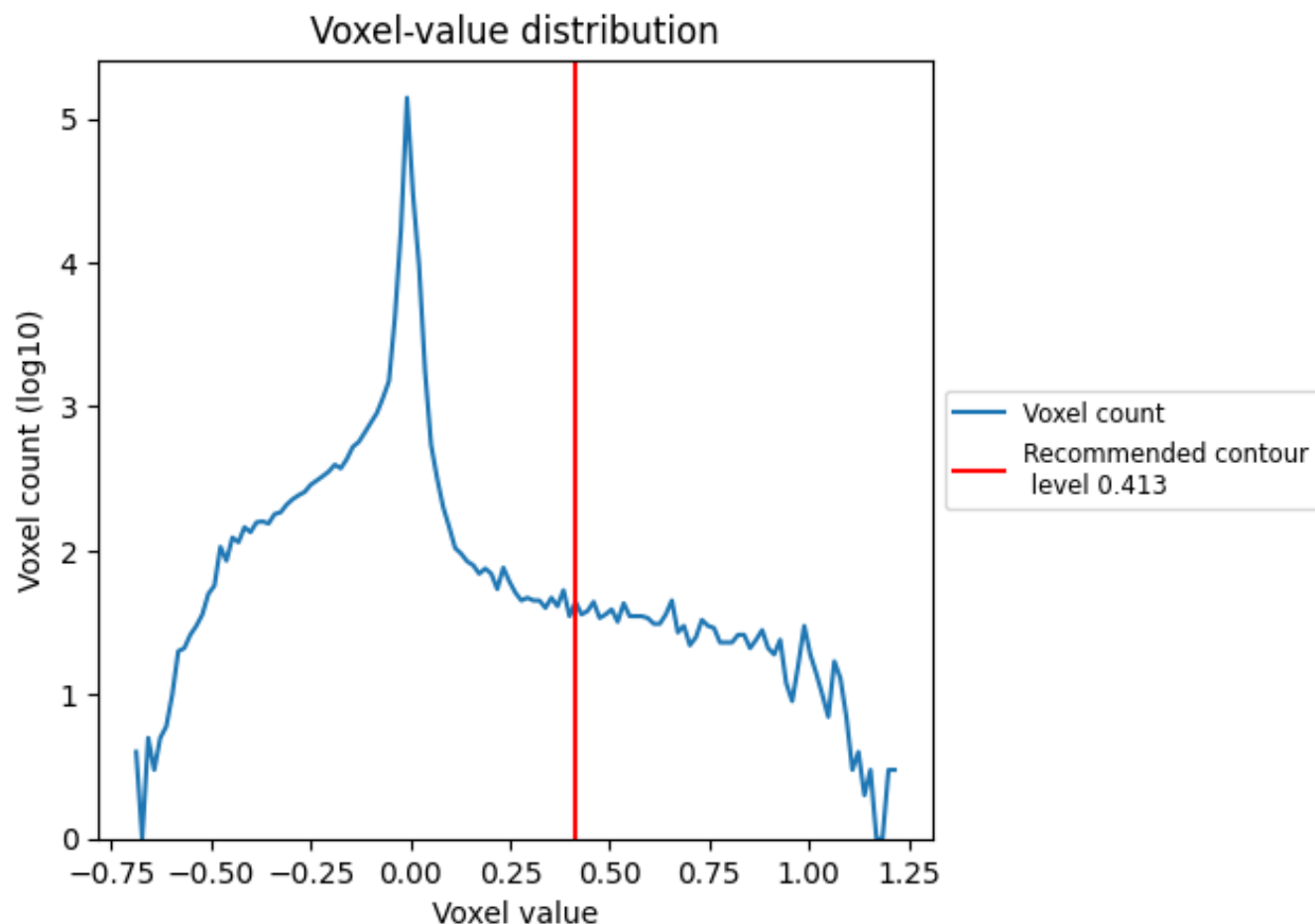
6.5 Mask visualisation [i](#)

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

7 Tomogram analysis [i](#)

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

7.1 Voxel-value distribution [i](#)



The voxel-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic.

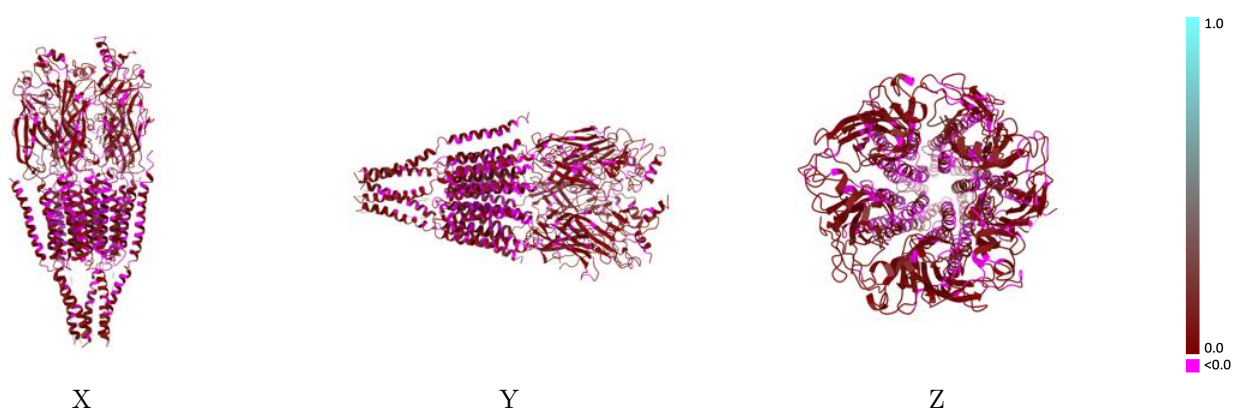
8 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-2377 and PDB model 4BOI. Per-residue inclusion information can be found in section 3 on page 4.

8.1 Map-model overlay [i](#)

This section was not generated.

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

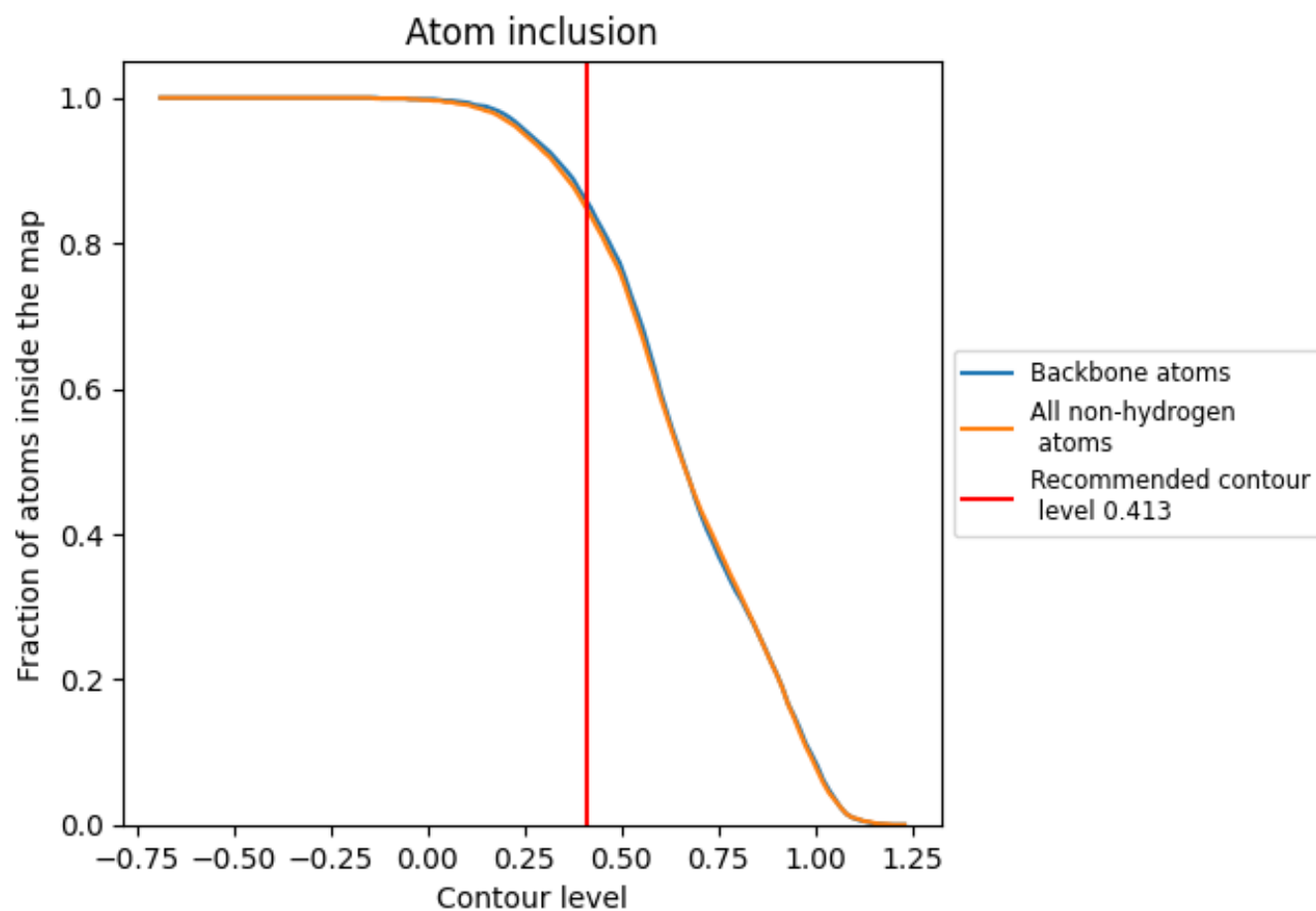


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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

This section was not generated.

8.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 84% of all non-hydrogen atoms, are inside the map.

8.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.413) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8450	0.0430
A	0.7690	0.0400
B	0.8790	0.0430
C	0.8520	0.0460
D	0.8310	0.0430
E	0.8950	0.0450

