



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 11:33 AM UTC

PDB ID : 5IPU / pdb_00005ipu
EMDB ID : EMD-8105
Title : Cryo-EM structure of GluN1/GluN2B NMDA receptor in the DCKA/D-APV-bound conformation, state 6
Authors : Zhu, S.; Stein, A.R.; Yoshioka, C.; Lee, C.H.; Goehring, A.; Mchaourab, S.H.; Gouaux, E.
Deposited on : 2016-03-09
Resolution : 15.40 Å(reported)

This is a Full wwPDB EM Validation 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:

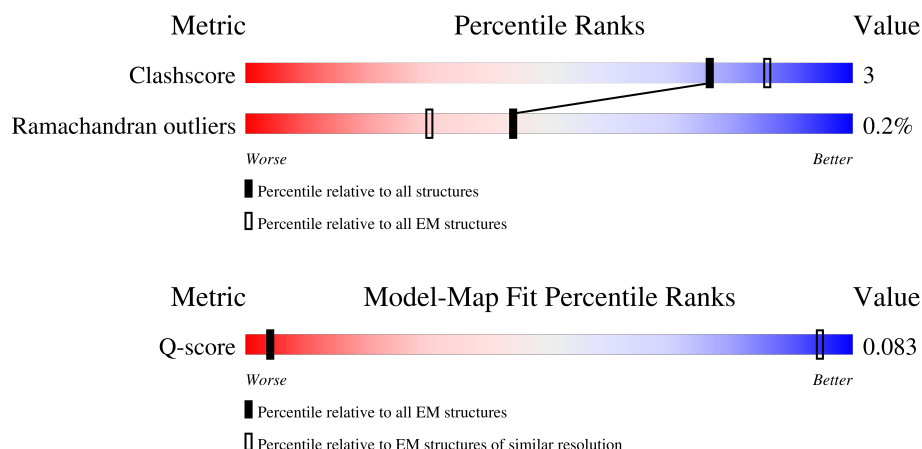
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 15.40 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	-
Q-score	-	25397	36 (14.90 - 15.90)

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	822	6% 76% 21%
1	C	822	7% 72% 24%
2	B	825	10% 72% 23%
2	D	825	12% 71% 24%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10156 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 N-methyl-D-aspartate receptor subunit NR1-8a.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	653	Total	C	N	O	0	0
			2612	1306	653	653		
1	C	626	Total	C	N	O	0	0
			2504	1252	626	626		

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	PHE	LYS	engineered mutation	UNP C0KD18
A	52	PHE	ARG	engineered mutation	UNP C0KD18
A	300	GLN	ASN	engineered mutation	UNP C0KD18
A	350	GLN	ASN	engineered mutation	UNP C0KD18
A	368	ASP	ASN	engineered mutation	UNP C0KD18
A	440	ASP	ASN	engineered mutation	UNP C0KD18
A	469	ASP	ASN	engineered mutation	UNP C0KD18
A	493	ALA	LYS	engineered mutation	UNP C0KD18
A	494	ALA	LYS	engineered mutation	UNP C0KD18
A	495	ALA	GLU	engineered mutation	UNP C0KD18
A	?	-	LYS	deletion	UNP C0KD18
A	?	-	VAL	deletion	UNP C0KD18
A	?	-	ASN	deletion	UNP C0KD18
A	?	-	SER	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	?	-	GLU	deletion	UNP C0KD18
A	602	ARG	GLY	engineered mutation	UNP C0KD18
A	609	LEU	ILE	engineered mutation	UNP C0KD18
A	648	ARG	ASP	engineered mutation	UNP C0KD18
A	761	GLU	ASN	engineered mutation	UNP C0KD18
A	829	SER	-	expression tag	UNP C0KD18
A	830	ARG	-	expression tag	UNP C0KD18
A	831	ALA	-	expression tag	UNP C0KD18
A	832	GLU	-	expression tag	UNP C0KD18

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Chain	Residue	Modelled	Actual	Comment	Reference
A	833	ALA	-	expression tag	UNP C0KD18
A	834	LYS	-	expression tag	UNP C0KD18
A	835	ARG	-	expression tag	UNP C0KD18
A	836	MET	-	expression tag	UNP C0KD18
A	837	LYS	-	expression tag	UNP C0KD18
A	838	GLY	-	expression tag	UNP C0KD18
A	839	LEU	-	expression tag	UNP C0KD18
A	840	GLU	-	expression tag	UNP C0KD18
A	841	VAL	-	expression tag	UNP C0KD18
A	842	LEU	-	expression tag	UNP C0KD18
A	843	PHE	-	expression tag	UNP C0KD18
A	844	GLN	-	expression tag	UNP C0KD18
C	51	PHE	LYS	engineered mutation	UNP C0KD18
C	52	PHE	ARG	engineered mutation	UNP C0KD18
C	300	GLN	ASN	engineered mutation	UNP C0KD18
C	350	GLN	ASN	engineered mutation	UNP C0KD18
C	368	ASP	ASN	engineered mutation	UNP C0KD18
C	440	ASP	ASN	engineered mutation	UNP C0KD18
C	469	ASP	ASN	engineered mutation	UNP C0KD18
C	493	ALA	LYS	engineered mutation	UNP C0KD18
C	494	ALA	LYS	engineered mutation	UNP C0KD18
C	495	ALA	GLU	engineered mutation	UNP C0KD18
C	?	-	LYS	deletion	UNP C0KD18
C	?	-	VAL	deletion	UNP C0KD18
C	?	-	ASN	deletion	UNP C0KD18
C	?	-	SER	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	?	-	GLU	deletion	UNP C0KD18
C	602	ARG	GLY	engineered mutation	UNP C0KD18
C	609	LEU	ILE	engineered mutation	UNP C0KD18
C	648	ARG	ASP	engineered mutation	UNP C0KD18
C	761	GLU	ASN	engineered mutation	UNP C0KD18
C	829	SER	-	expression tag	UNP C0KD18
C	830	ARG	-	expression tag	UNP C0KD18
C	831	ALA	-	expression tag	UNP C0KD18
C	832	GLU	-	expression tag	UNP C0KD18
C	833	ALA	-	expression tag	UNP C0KD18
C	834	LYS	-	expression tag	UNP C0KD18
C	835	ARG	-	expression tag	UNP C0KD18
C	836	MET	-	expression tag	UNP C0KD18

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Chain	Residue	Modelled	Actual	Comment	Reference
C	837	LYS	-	expression tag	UNP C0KD18
C	838	GLY	-	expression tag	UNP C0KD18
C	839	LEU	-	expression tag	UNP C0KD18
C	840	GLU	-	expression tag	UNP C0KD18
C	841	VAL	-	expression tag	UNP C0KD18
C	842	LEU	-	expression tag	UNP C0KD18
C	843	PHE	-	expression tag	UNP C0KD18
C	844	GLN	-	expression tag	UNP C0KD18

- Molecule 2 is a protein called Ionotropic glutamate receptor subunit NR2B.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	B	633	Total	C	N	O	0	0
			2532	1266	633	633		
2	D	627	Total	C	N	O	0	0
			2508	1254	627	627		

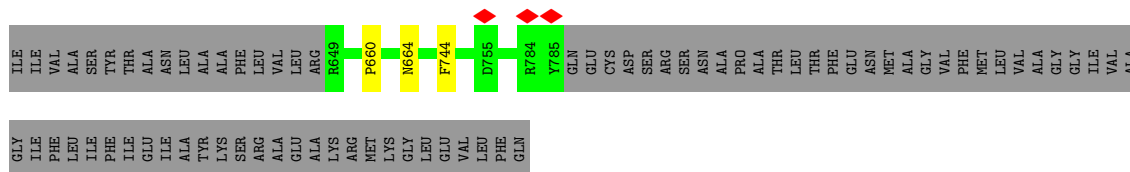
There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	20	SER	MET	engineered mutation	UNP A7XY94
B	21	ARG	GLY	engineered mutation	UNP A7XY94
B	22	ALA	CYS	engineered mutation	UNP A7XY94
B	64	GLU	ALA	engineered mutation	UNP A7XY94
B	69	GLN	ASN	engineered mutation	UNP A7XY94
B	343	ASP	ASN	engineered mutation	UNP A7XY94
B	?	-	LYS	deletion	UNP A7XY94
B	?	-	TYR	deletion	UNP A7XY94
B	?	-	TYR	deletion	UNP A7XY94
B	?	-	VAL	deletion	UNP A7XY94
B	486	VAL	THR	engineered mutation	UNP A7XY94
B	?	-	ARG	deletion	UNP A7XY94
B	?	-	CYS	deletion	UNP A7XY94
B	?	-	LEU	deletion	UNP A7XY94
B	?	-	ALA	deletion	UNP A7XY94
B	?	-	ASP	deletion	UNP A7XY94
B	?	-	GLY	deletion	UNP A7XY94
B	?	-	ARG	deletion	UNP A7XY94
B	?	-	GLU	deletion	UNP A7XY94
B	?	-	PRO	deletion	UNP A7XY94
B	?	-	GLY	deletion	UNP A7XY94
B	601	LEU	VAL	engineered mutation	UNP A7XY94

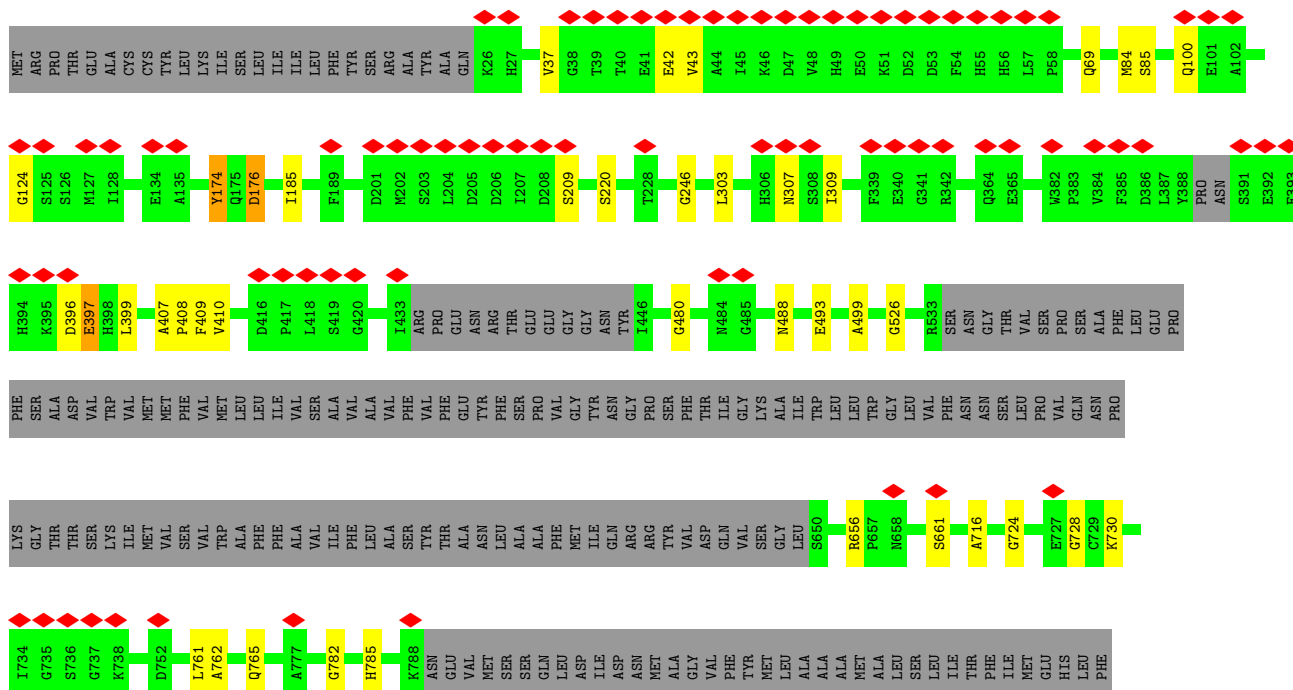
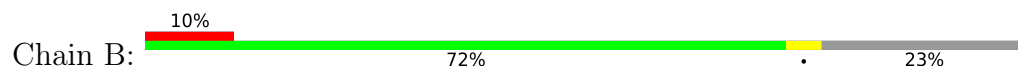
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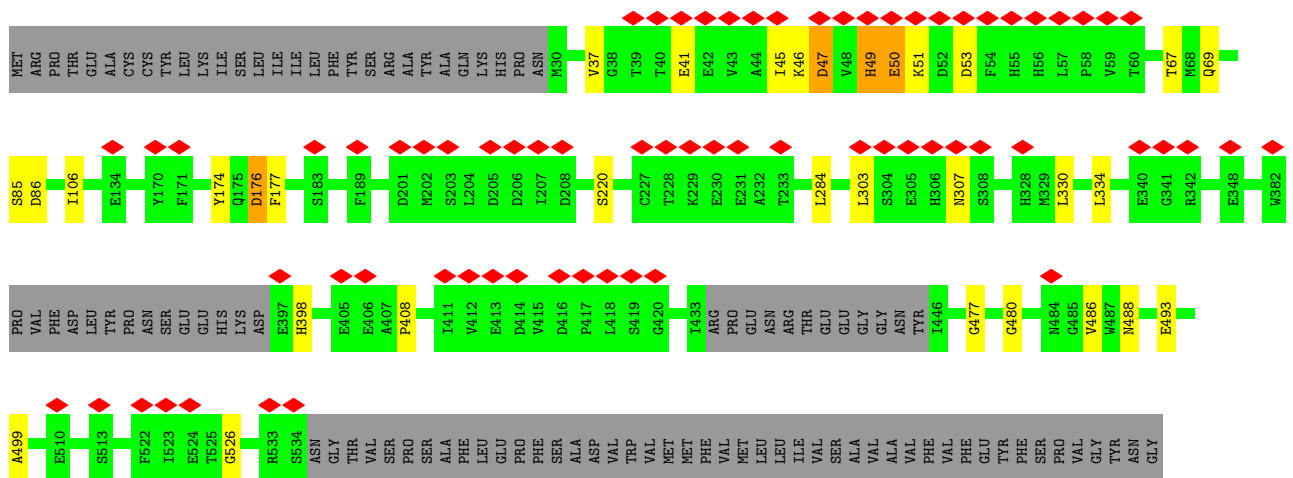
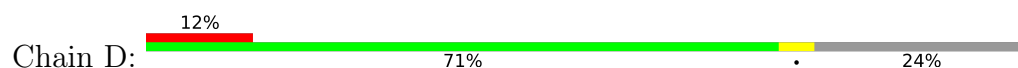
Chain	Residue	Modelled	Actual	Comment	Reference
B	640	ARG	GLU	engineered mutation	UNP A7XY94
B	641	ARG	GLU	engineered mutation	UNP A7XY94
D	20	SER	MET	engineered mutation	UNP A7XY94
D	21	ARG	GLY	engineered mutation	UNP A7XY94
D	22	ALA	CYS	engineered mutation	UNP A7XY94
D	64	GLU	ALA	engineered mutation	UNP A7XY94
D	69	GLN	ASN	engineered mutation	UNP A7XY94
D	343	ASP	ASN	engineered mutation	UNP A7XY94
D	?	-	LYS	deletion	UNP A7XY94
D	?	-	TYR	deletion	UNP A7XY94
D	?	-	TYR	deletion	UNP A7XY94
D	?	-	VAL	deletion	UNP A7XY94
D	486	VAL	THR	engineered mutation	UNP A7XY94
D	?	-	ARG	deletion	UNP A7XY94
D	?	-	CYS	deletion	UNP A7XY94
D	?	-	LEU	deletion	UNP A7XY94
D	?	-	ALA	deletion	UNP A7XY94
D	?	-	ASP	deletion	UNP A7XY94
D	?	-	GLY	deletion	UNP A7XY94
D	?	-	ARG	deletion	UNP A7XY94
D	?	-	GLU	deletion	UNP A7XY94
D	?	-	PRO	deletion	UNP A7XY94
D	?	-	GLY	deletion	UNP A7XY94
D	601	LEU	VAL	engineered mutation	UNP A7XY94
D	640	ARG	GLU	engineered mutation	UNP A7XY94
D	641	ARG	GLU	engineered mutation	UNP A7XY94

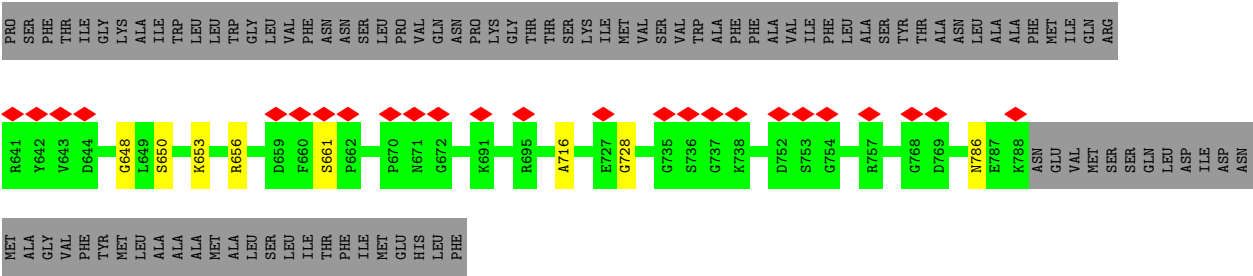


• Molecule 2: Ionotropic glutamate receptor subunit NR2B



• Molecule 2: Ionotropic glutamate receptor subunit NR2B





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	18309	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 ($e^-/\text{\AA}^2$)	10.5	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.139	Depositor
Minimum map value	-0.041	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.052	Depositor
Map size ($\text{\AA}$)	236.47003, 215.07004, 204.37003	wwPDB
Map dimensions	221, 201, 191	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.0700002, 1.0700002, 1.0700002	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	0.52	0/2608	1.29	20/3253 (0.6%)
1	C	0.53	0/2500	1.38	21/3118 (0.7%)
2	B	0.51	0/2528	1.46	27/3153 (0.9%)
2	D	0.54	1/2504 (0.0%)	1.41	27/3123 (0.9%)
All	All	0.52	1/10140 (0.0%)	1.39	95/12647 (0.8%)

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
2	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	49	HIS	CA-C	-5.42	1.46	1.52

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	409	PHE	N-CA-C	-26.30	71.94	108.23
2	B	410	VAL	N-CA-C	12.90	136.17	109.34
2	D	176	ASP	N-CA-C	-12.87	98.05	113.88
1	C	103	THR	CA-C-N	-12.73	104.72	121.91
1	C	103	THR	C-N-CA	-12.73	104.72	121.91
2	B	176	ASP	N-CA-C	-11.65	98.56	113.12
2	B	69	GLN	N-CA-C	10.55	122.86	111.36
2	D	69	GLN	N-CA-C	-8.99	101.56	111.36
1	A	381	TRP	CA-C-N	-8.86	110.91	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	381	TRP	C-N-CA	-8.86	110.91	119.85
2	D	41	GLU	N-CA-C	7.92	122.52	109.06
2	B	42	GLU	N-CA-C	-7.76	96.92	109.96
2	D	656	ARG	CA-C-N	7.33	126.84	118.85
2	D	656	ARG	C-N-CA	7.33	126.84	118.85
2	D	661	SER	CA-C-N	-7.26	112.52	119.85
2	D	661	SER	C-N-CA	-7.26	112.52	119.85
1	C	123	THR	N-CA-C	7.19	126.12	110.80
2	B	174	TYR	N-CA-C	-6.74	101.00	110.50
2	B	656	ARG	CA-C-N	6.69	126.14	118.85
2	B	656	ARG	C-N-CA	6.69	126.14	118.85
2	B	37	VAL	CA-C-N	-6.67	108.33	121.41
2	B	37	VAL	C-N-CA	-6.67	108.33	121.41
1	C	412	ARG	CA-C-N	-6.65	113.09	120.14
1	C	412	ARG	C-N-CA	-6.65	113.09	120.14
1	A	159	ASN	CA-C-N	-6.62	113.96	122.77
1	A	159	ASN	C-N-CA	-6.62	113.96	122.77
1	C	398	LEU	N-CA-C	6.62	119.68	108.90
2	D	86	ASP	N-CA-C	6.60	121.42	113.23
1	A	192	ASP	N-CA-C	-6.51	104.11	111.14
2	B	396	ASP	CA-C-N	6.48	133.93	121.54
2	B	396	ASP	C-N-CA	6.48	133.93	121.54
1	C	446	ARG	CA-C-N	6.47	126.99	120.14
1	C	446	ARG	C-N-CA	6.47	126.99	120.14
1	C	521	ARG	N-CA-C	6.46	118.40	111.36
2	B	785	HIS	N-CA-C	-6.28	104.20	112.24
2	D	728	GLY	N-CA-C	-6.14	101.81	110.96
2	D	177	PHE	N-CA-C	-6.14	104.51	111.14
1	C	379	ILE	CA-C-N	-6.10	114.64	123.06
1	C	379	ILE	C-N-CA	-6.10	114.64	123.06
1	A	657	ILE	N-CA-C	-6.08	104.42	110.62
2	D	46	LYS	N-CA-C	-6.08	104.22	111.69
1	C	223	ALA	N-CA-C	6.07	117.46	108.60
1	C	23	ASP	CA-C-N	6.01	126.20	120.31
1	C	23	ASP	C-N-CA	6.01	126.20	120.31
2	D	486	VAL	N-CA-C	6.01	117.25	108.53
2	B	661	SER	CA-C-N	-5.99	113.80	119.85
2	B	661	SER	C-N-CA	-5.99	113.80	119.85
2	D	85	SER	N-CA-C	5.99	118.57	111.33
1	A	446	ARG	CA-C-N	5.97	126.47	120.14
1	A	446	ARG	C-N-CA	5.97	126.47	120.14
2	B	185	ILE	N-CA-C	5.84	116.62	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	399	LEU	N-CA-C	5.81	118.19	108.02
2	D	786	ASN	N-CA-C	-5.81	104.55	111.69
2	D	50	GLU	N-CA-C	-5.80	105.69	112.89
1	A	23	ASP	CA-C-N	-5.79	114.63	120.31
1	A	23	ASP	C-N-CA	-5.79	114.63	120.31
2	B	396	ASP	N-CA-C	5.76	116.98	108.86
2	D	284	LEU	N-CA-C	-5.71	105.04	113.16
1	A	744	PHE	N-CA-C	-5.70	104.28	111.92
1	A	278	SER	N-CA-C	5.62	117.08	111.07
1	A	756	SER	CA-C-N	5.62	125.71	119.87
1	A	756	SER	C-N-CA	5.62	125.71	119.87
2	D	47	ASP	N-CA-C	-5.54	105.24	111.28
2	D	477	GLY	N-CA-C	5.52	123.24	115.30
1	C	156	ARG	N-CA-C	-5.49	105.29	111.28
1	C	151	TRP	N-CA-C	-5.46	105.32	111.28
2	D	398	HIS	N-CA-C	-5.46	104.81	112.12
1	C	328	SER	N-CA-C	5.45	119.64	113.15
1	A	188	GLU	N-CA-C	-5.40	102.40	110.23
2	B	397	GLU	N-CA-C	5.39	122.28	110.80
1	C	204	LEU	N-CA-C	5.34	119.76	112.88
1	A	198	GLU	N-CA-C	-5.31	102.81	110.40
1	C	108	SER	N-CA-C	-5.30	105.51	111.28
2	B	309	ILE	CA-C-N	5.27	124.88	119.56
2	B	309	ILE	C-N-CA	5.27	124.88	119.56
2	D	653	LYS	N-CA-C	-5.27	105.54	111.28
2	D	106	ILE	N-CA-C	-5.25	105.38	110.42
2	B	785	HIS	CA-C-N	5.19	128.19	120.31
2	B	785	HIS	C-N-CA	5.19	128.19	120.31
1	A	397	ARG	N-CA-C	5.16	118.15	109.95
1	C	132	SER	N-CA-C	5.16	117.30	111.11
2	D	37	VAL	CA-C-N	5.15	133.06	120.28
2	D	37	VAL	C-N-CA	5.15	133.06	120.28
1	A	486	GLU	N-CA-C	5.13	115.01	108.24
1	A	137	PHE	N-CA-C	5.09	117.06	108.96
2	B	765	GLN	N-CA-C	-5.09	105.82	112.23
1	C	744	PHE	N-CA-C	-5.09	105.10	111.92
1	A	255	SER	N-CA-C	5.08	117.41	110.55
2	B	209	SER	N-CA-C	-5.06	104.66	111.54
2	B	220	SER	CA-C-N	5.05	124.54	119.19
2	B	220	SER	C-N-CA	5.05	124.54	119.19
2	D	67	THR	CA-C-N	-5.01	115.02	122.74
2	D	67	THR	C-N-CA	-5.01	115.02	122.74

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	220	SER	CA-C-N	5.01	124.50	119.19
2	D	220	SER	C-N-CA	5.01	124.50	119.19

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	542	LYS	Peptide
2	B	246	GLY	Peptide

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	2612	0	699	8	0
1	C	2504	0	664	8	0
2	B	2532	0	687	10	0
2	D	2508	0	685	12	0
All	All	10156	0	2735	38	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.

All (38) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance ($\text{\AA}$)	Clash overlap ($\text{\AA}$)
1:C:660:PRO:O	1:C:664:ASN:N	2.15	0.77
2:D:50:GLU:O	2:D:53:ASP:N	2.20	0.75
1:A:660:PRO:O	1:A:664:ASN:N	2.22	0.71
1:C:464:LEU:O	1:C:468:MET:N	2.20	0.71
1:A:464:LEU:O	1:A:468:MET:N	2.23	0.71
1:A:483:GLY:HA2	1:A:497:ASN:O	1.93	0.69
2:D:47:ASP:O	2:D:51:LYS:N	2.29	0.66
2:D:330:LEU:O	2:D:334:LEU:N	2.30	0.63
1:C:265:GLY:HA3	1:C:381:TRP:C	2.23	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:493:GLU:O	2:D:499:ALA:N	2.31	0.63
2:D:480:GLY:HA2	2:D:488:ASN:O	1.98	0.63
2:D:45:ILE:O	2:D:49:HIS:N	2.32	0.62
2:D:526:GLY:O	2:D:716:ALA:N	2.36	0.58
2:B:303:LEU:O	2:B:307:ASN:N	2.34	0.57
2:B:526:GLY:O	2:B:716:ALA:N	2.37	0.57
1:C:102:LEU:O	1:C:105:THR:N	2.37	0.57
2:B:493:GLU:O	2:B:499:ALA:N	2.37	0.57
1:A:498:GLY:O	1:A:501:GLY:N	2.38	0.56
1:C:149:LEU:O	1:C:153:GLU:N	2.35	0.55
1:C:150:VAL:O	1:C:154:MET:N	2.35	0.55
1:A:365:GLY:HA2	1:A:375:ASN:H	1.73	0.54
1:C:365:GLY:HA2	1:C:375:ASN:H	1.75	0.52
2:D:303:LEU:O	2:D:307:ASN:N	2.36	0.52
1:A:30:GLY:HA2	1:A:63:THR:O	2.11	0.51
2:D:174:TYR:C	2:D:176:ASP:H	2.17	0.51
2:B:480:GLY:HA2	2:B:488:ASN:O	2.12	0.49
2:B:724:GLY:HA3	2:B:782:GLY:HA3	1.94	0.49
2:B:728:GLY:C	2:B:730:LYS:H	2.21	0.49
2:B:174:TYR:C	2:B:176:ASP:H	2.24	0.46
2:D:648:GLY:C	2:D:650:SER:H	2.25	0.44
1:A:709:GLN:O	1:A:713:ASP:N	2.41	0.43
2:B:84:MET:O	2:B:85:SER:C	2.63	0.42
2:D:50:GLU:O	2:D:51:LYS:C	2.63	0.42
2:B:761:LEU:O	2:B:762:ALA:C	2.64	0.41
1:C:383:GLY:C	1:C:385:GLU:H	2.28	0.41
2:B:100:GLN:O	2:B:124:GLY:HA3	2.20	0.41
2:D:174:TYR:C	2:D:176:ASP:N	2.79	0.40
1:A:482:PHE:O	1:A:499:MET:N	2.53	0.40

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	645/822 (78%)	626 (97%)	19 (3%)	0	100	100
1	C	618/822 (75%)	597 (97%)	21 (3%)	0	100	100
2	B	625/825 (76%)	592 (95%)	29 (5%)	4 (1%)	21	59
2	D	619/825 (75%)	592 (96%)	26 (4%)	1 (0%)	43	78
All	All	2507/3294 (76%)	2407 (96%)	95 (4%)	5 (0%)	44	78

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	407	ALA
2	B	408	PRO
2	D	408	PRO
2	B	397	GLU
2	B	43	VAL

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 ⓘ

There are no chain breaks in this entry.

6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8105. 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](#)

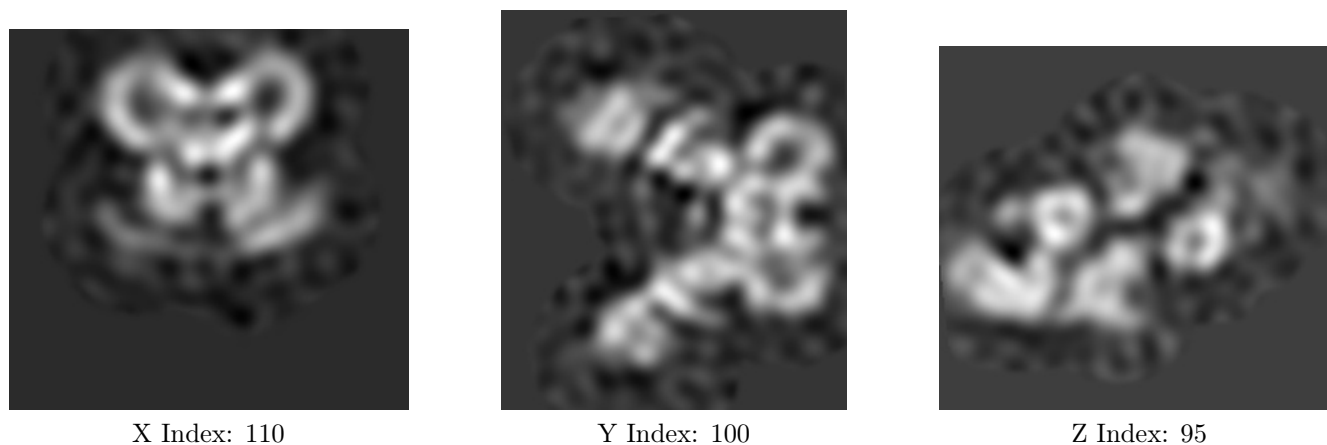
6.1.1 Primary map



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



X Index: 109



Y Index: 89

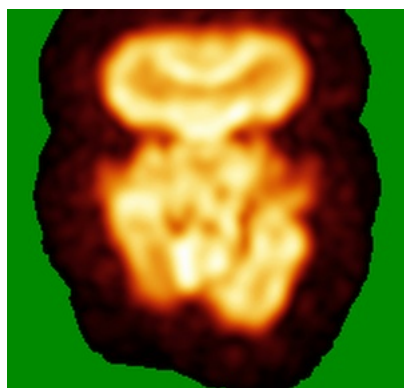


Z Index: 140

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

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

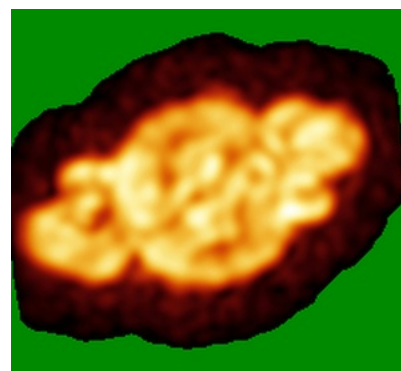
6.4.1 Primary map



X



Y

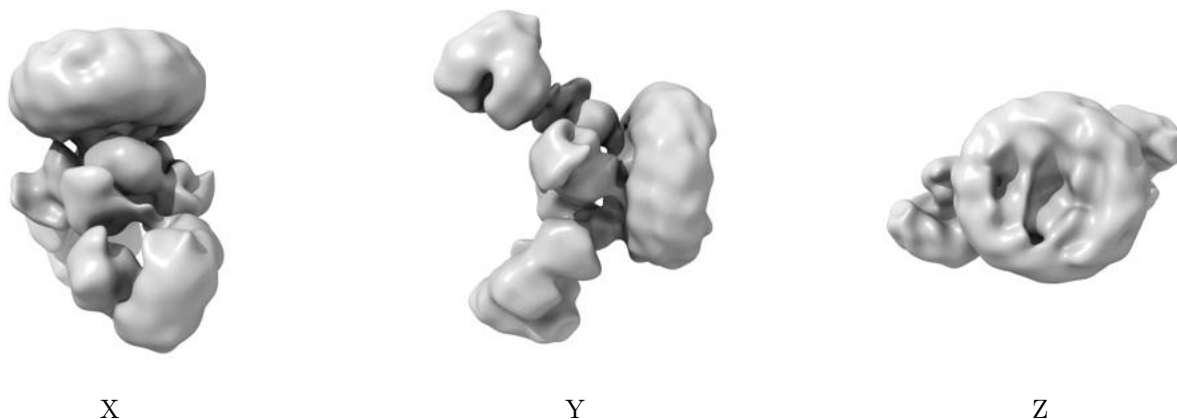


Z

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

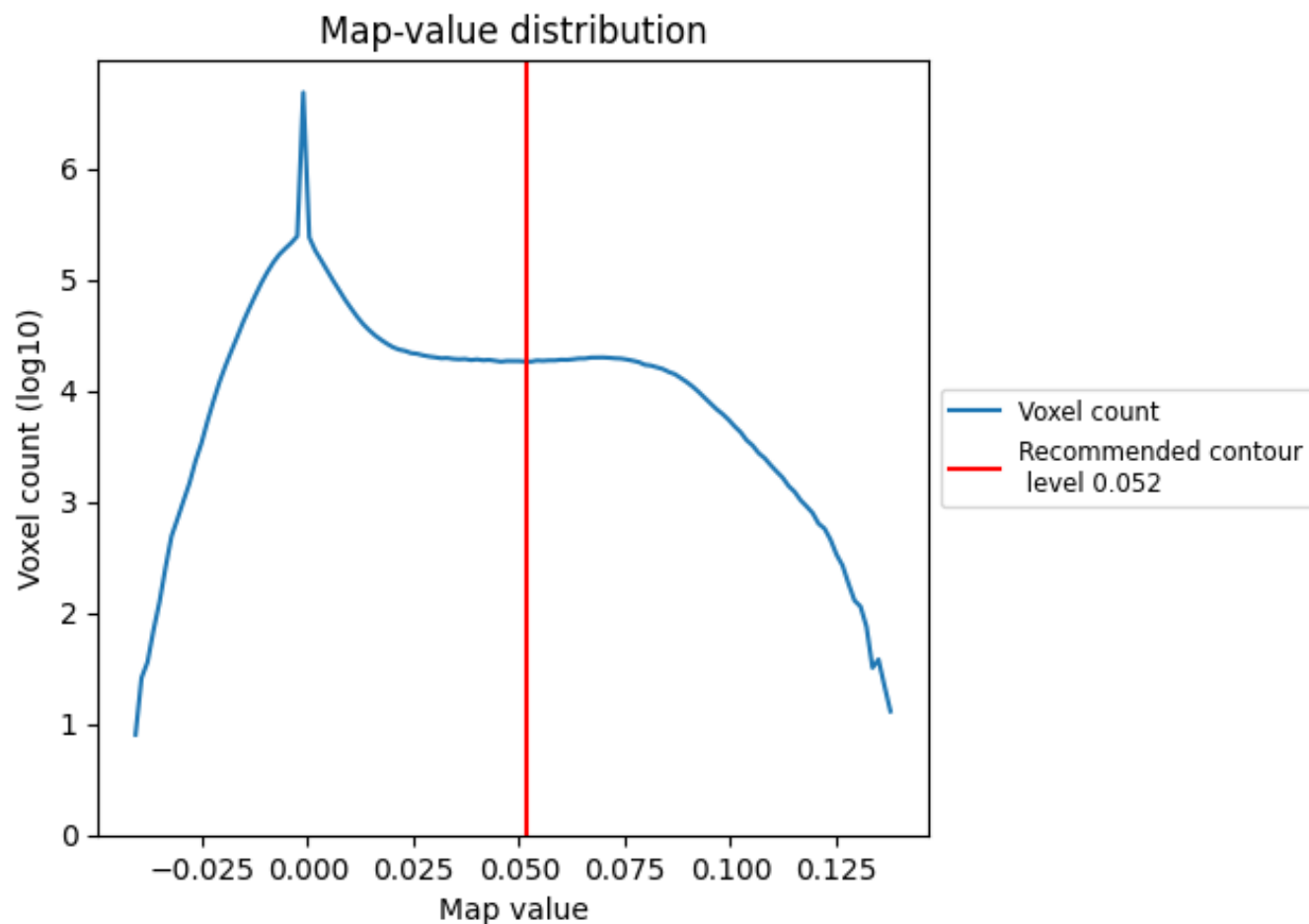
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

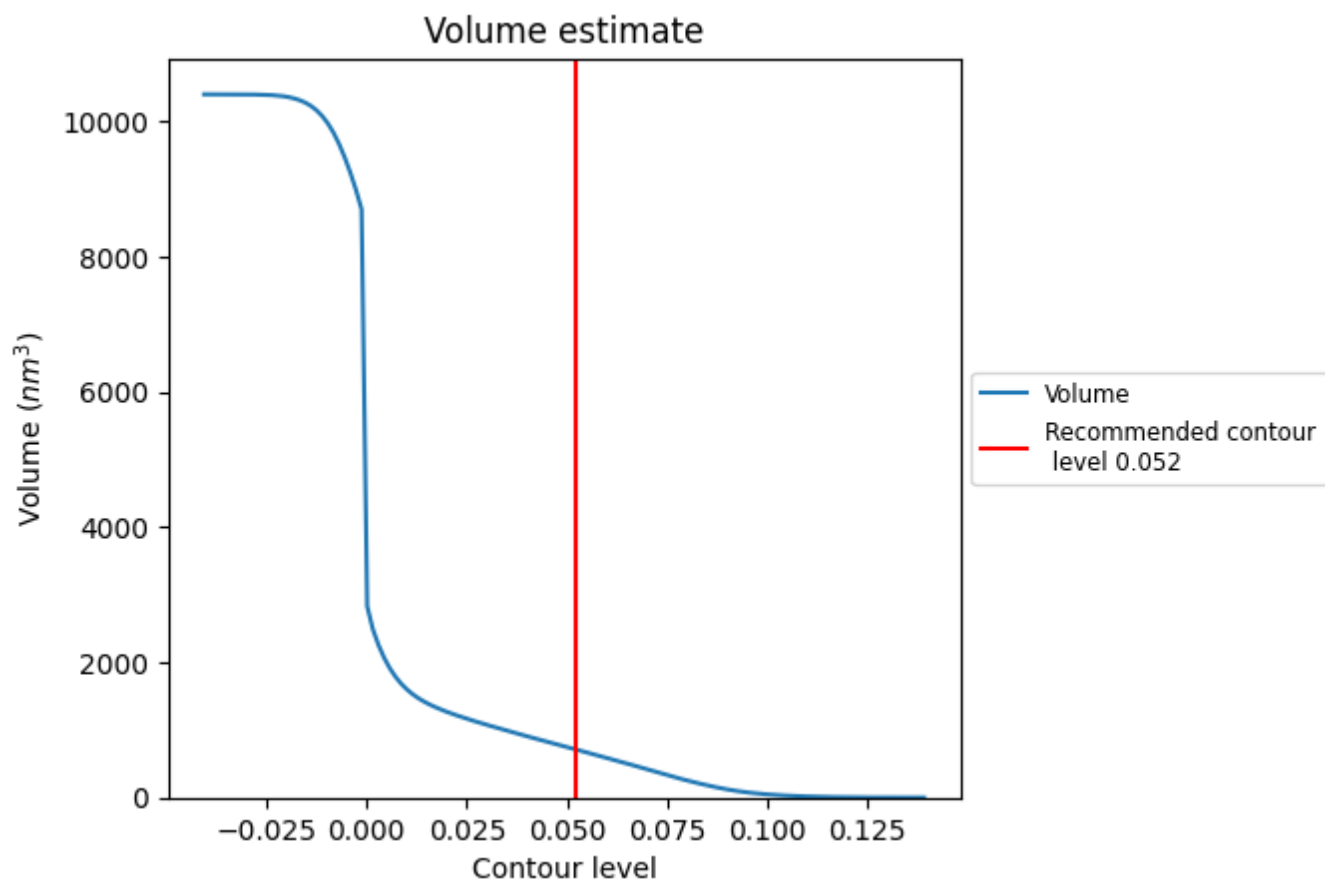
This section contains the results of statistical analysis of the map.

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 714 nm³; this corresponds to an approximate mass of 645 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](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

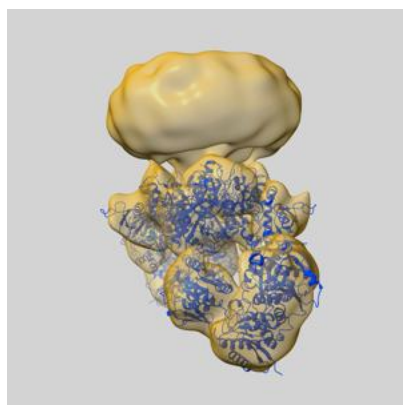
8 Fourier-Shell correlation ⓘ

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

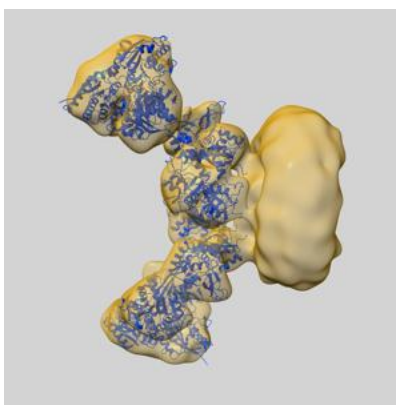
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8105 and PDB model 5IPU. Per-residue inclusion information can be found in section [3](#) on page [7](#).

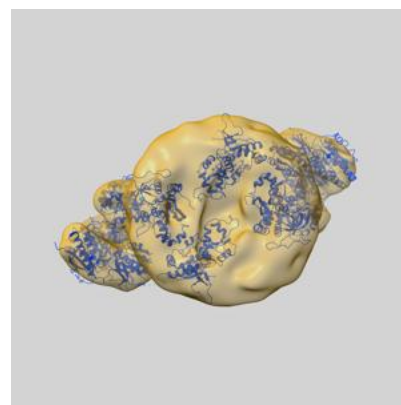
9.1 Map-model overlay [i](#)



X



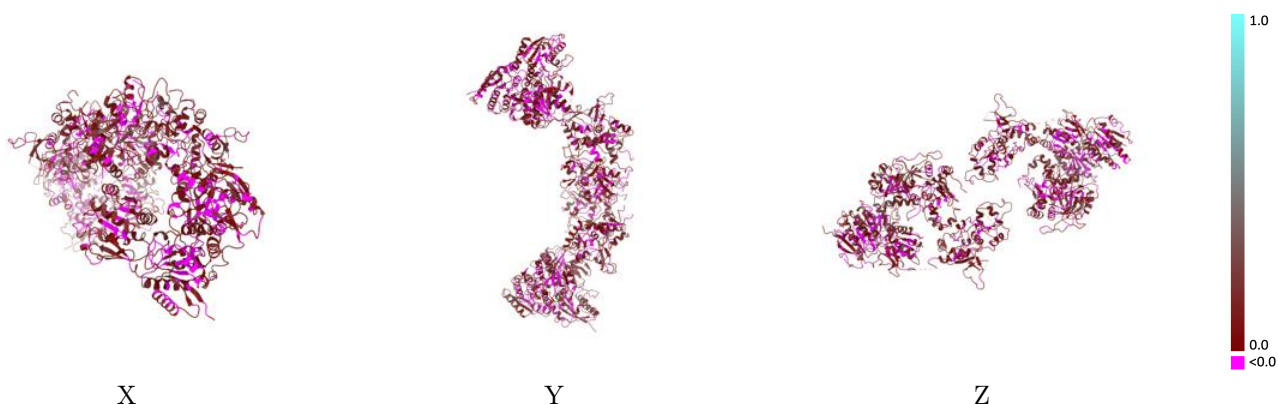
Y



Z

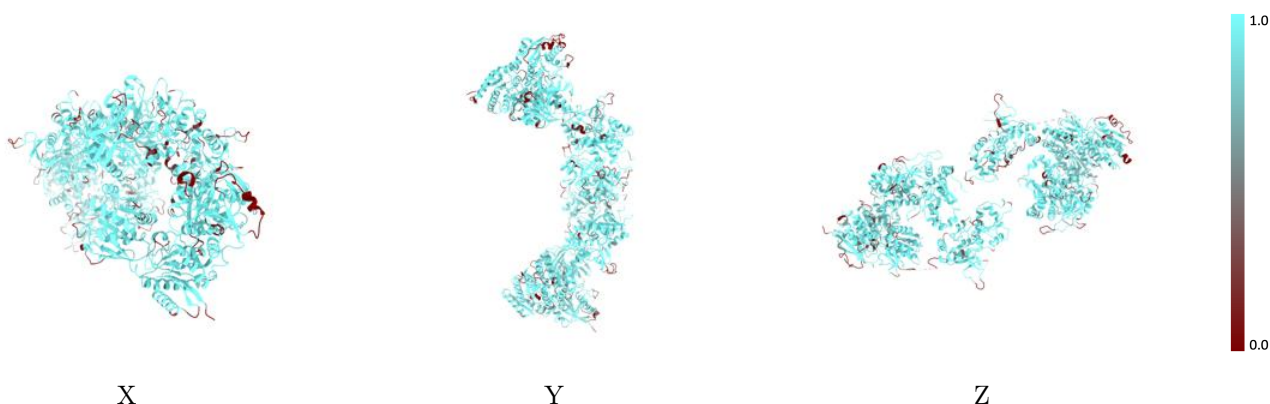
The images above show the 3D surface view of the map at the recommended contour level 0.052 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



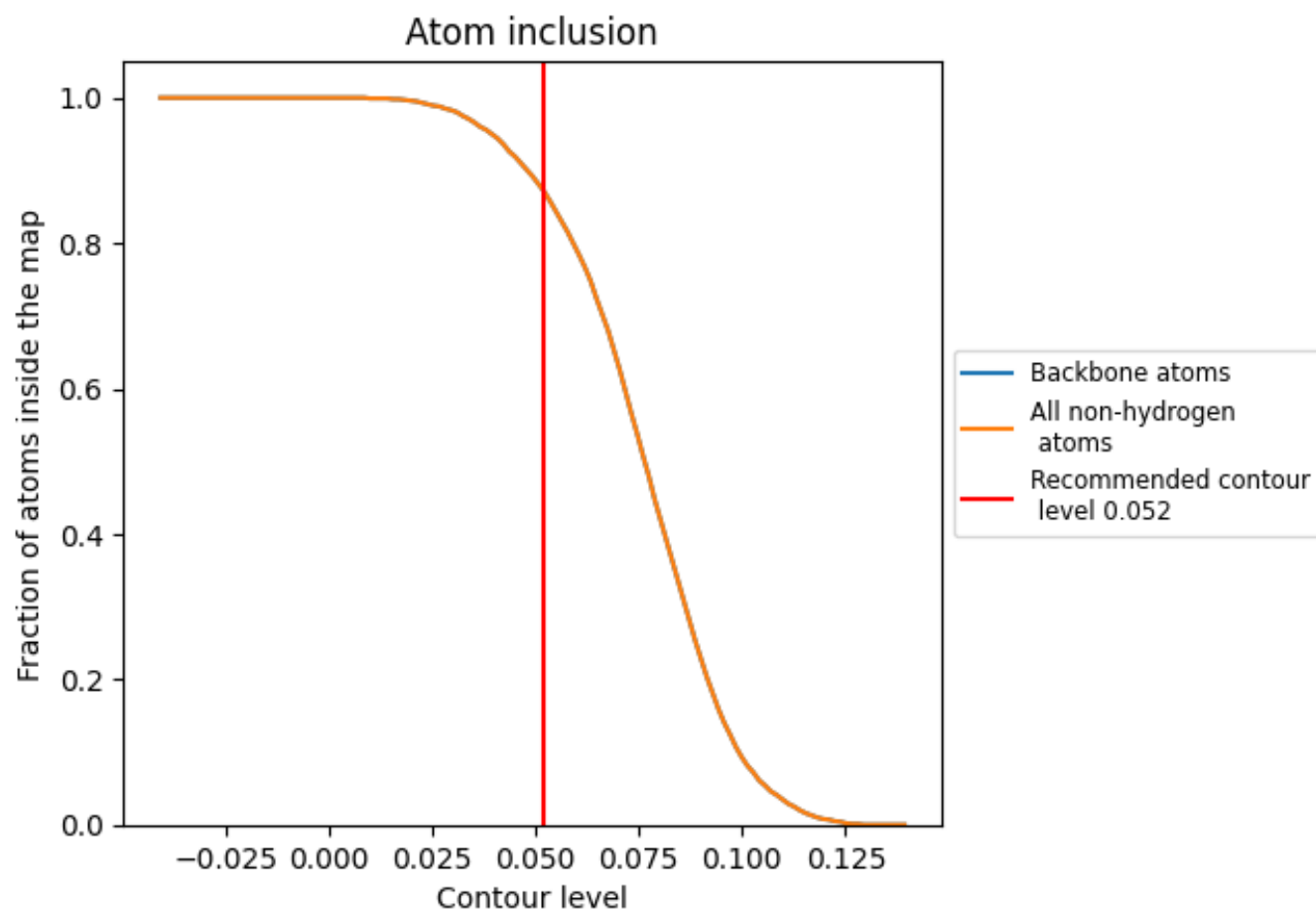
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.052).

9.4 Atom inclusion ⓘ



At the recommended contour level, 87% of all backbone atoms, 87% 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 (0.052) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8720	0.0830
A	0.9100	0.0920
B	0.8590	0.0830
C	0.8950	0.0830
D	0.8220	0.0750

