



## Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 09:57 PM UTC

PDB ID : 5IPS / pdb\_00005ips  
EMDB ID : EMD-8103  
Title : Cryo-EM structure of GluN1/GluN2B NMDA receptor in the DCKA/D-APV-bound conformation, state 4  
Authors : Zhu, S.; Stein, A.R.; Yoshioka, C.; Lee, C.H.; Goehring, A.; Mchaourab, S.H.; Gouaux, E.  
Deposited on : 2016-03-09  
Resolution : 13.50 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

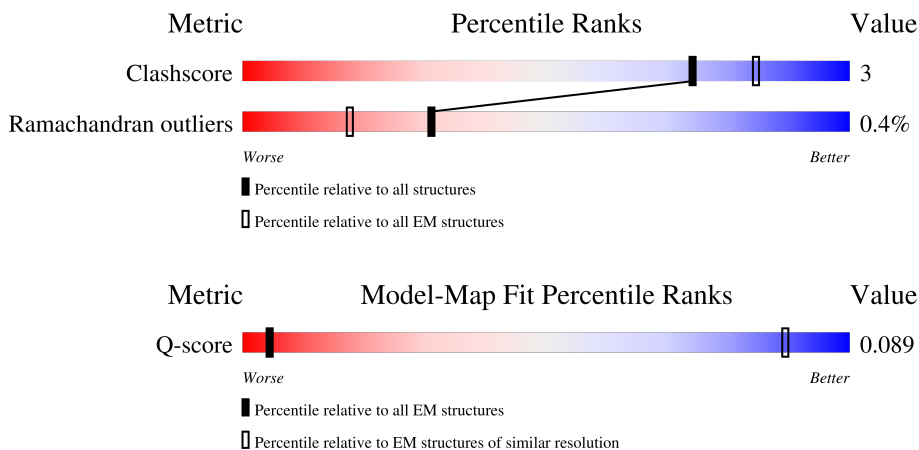
EMDB validation analysis : 0.0.1.dev132  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)  
MapQ : 1.9.13  
Ideal geometry (proteins)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

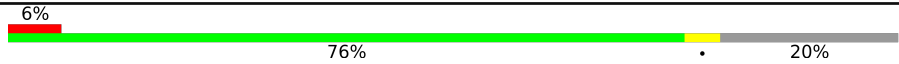



The reported resolution of this entry is 13.50 Å.

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	74 ( 13.00 - 14.00 )

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  $\geq 3$ , 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% 20%
1	C	822	 11% 72% 24%
2	B	825	 12% 72% 23%
2	D	825	 17% 71% 5% 24%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10181 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
			Total	C	N	O		
1	A	658	2632	1316	658	658	0	0
1	C	627	2508	1254	627	627	0	0

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	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
			Total	C	N	O		
2	B	633	2532	1266	633	633	0	0
2	D	627	2509	1255	627	627	0	0

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







## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	36878	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	0.205	Depositor
Minimum map value	-0.085	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.022	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 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	0.52	0/2629	1.28	19/3281 (0.6%)
1	C	0.52	0/2503	1.38	21/3120 (0.7%)
2	B	0.53	1/2528 (0.0%)	1.56	28/3153 (0.9%)
2	D	0.54	1/2505 (0.0%)	1.41	28/3125 (0.9%)
All	All	0.53	2/10165 (0.0%)	1.41	96/12679 (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	2
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	409	PHE	C-N	-5.65	1.25	1.33
2	D	49	HIS	CA-C	-5.45	1.46	1.52

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	409	PHE	N-CA-C	27.54	149.30	112.68
2	B	409	PHE	CA-C-N	23.23	163.79	121.97
2	B	409	PHE	C-N-CA	23.23	163.79	121.97
1	C	103	THR	CA-C-N	-12.70	104.76	121.91
1	C	103	THR	C-N-CA	-12.70	104.76	121.91
2	B	69	GLN	N-CA-C	10.62	122.94	111.36
2	D	176	ASP	N-CA-C	-10.45	100.67	113.41
2	B	176	ASP	N-CA-C	-10.22	100.34	113.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	69	GLN	N-CA-C	-8.97	101.58	111.36
1	A	381	TRP	CA-C-N	-8.84	110.93	119.85
1	A	381	TRP	C-N-CA	-8.84	110.93	119.85
1	C	123	THR	N-CA-C	8.24	128.36	110.80
2	D	41	GLU	N-CA-C	7.91	122.50	109.06
2	B	42	GLU	N-CA-C	-7.76	96.92	109.96
2	D	656	ARG	CA-C-N	7.37	126.88	118.85
2	D	656	ARG	C-N-CA	7.37	126.88	118.85
2	D	661	SER	CA-C-N	-7.27	112.50	119.85
2	D	661	SER	C-N-CA	-7.27	112.50	119.85
2	B	174	TYR	N-CA-C	-6.78	100.93	110.50
2	B	656	ARG	CA-C-N	6.70	126.16	118.85
2	B	656	ARG	C-N-CA	6.70	126.16	118.85
2	B	37	VAL	CA-C-N	-6.68	108.32	121.41
2	B	37	VAL	C-N-CA	-6.68	108.32	121.41
1	C	412	ARG	CA-C-N	-6.64	113.10	120.14
1	C	412	ARG	C-N-CA	-6.64	113.10	120.14
2	D	86	ASP	N-CA-C	6.59	121.40	113.23
1	A	159	ASN	CA-C-N	-6.58	114.02	122.77
1	A	159	ASN	C-N-CA	-6.58	114.02	122.77
1	C	398	LEU	N-CA-C	6.57	119.61	108.90
1	A	192	ASP	N-CA-C	-6.54	104.08	111.14
1	C	446	ARG	CA-C-N	6.51	127.04	120.14
1	C	446	ARG	C-N-CA	6.51	127.04	120.14
2	B	396	ASP	CA-C-N	6.48	133.92	121.54
2	B	396	ASP	C-N-CA	6.48	133.92	121.54
1	C	521	ARG	N-CA-C	6.46	118.40	111.36
2	B	785	HIS	N-CA-C	-6.26	104.22	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
2	D	46	LYS	N-CA-C	-6.09	104.20	111.69
1	C	379	ILE	CA-C-N	-6.06	114.70	123.06
1	C	379	ILE	C-N-CA	-6.06	114.70	123.06
1	A	657	ILE	N-CA-C	-6.05	104.44	110.62
1	C	223	ALA	N-CA-C	6.03	117.41	108.60
2	D	85	SER	N-CA-C	6.03	118.62	111.33
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	B	661	SER	CA-C-N	-6.00	113.78	119.85
2	B	661	SER	C-N-CA	-6.00	113.78	119.85
2	D	486	VAL	N-CA-C	5.98	117.20	108.53
1	A	446	ARG	CA-C-N	5.96	126.45	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	446	ARG	C-N-CA	5.96	126.45	120.14
2	D	399	LEU	N-CA-C	5.90	123.38	110.80
2	D	786	ASN	N-CA-C	-5.83	104.51	111.69
2	D	50	GLU	N-CA-C	-5.83	105.67	112.89
2	B	185	ILE	N-CA-C	5.81	116.59	110.72
2	B	399	LEU	N-CA-C	5.80	118.16	108.02
2	B	396	ASP	N-CA-C	5.78	117.01	108.86
1	A	23	ASP	CA-C-N	-5.75	114.68	120.31
1	A	23	ASP	C-N-CA	-5.75	114.68	120.31
1	A	744	PHE	N-CA-C	-5.71	104.27	111.92
2	D	284	LEU	N-CA-C	-5.71	105.06	113.16
1	A	278	SER	N-CA-C	5.63	117.09	111.07
2	D	533	ARG	N-CA-C	5.60	122.72	110.80
2	D	47	ASP	N-CA-C	-5.57	105.20	111.28
2	D	477	GLY	N-CA-C	5.54	123.27	115.30
1	C	156	ARG	N-CA-C	-5.51	105.27	111.28
2	D	398	HIS	N-CA-C	-5.49	104.77	112.12
1	C	328	SER	N-CA-C	5.45	119.64	113.15
1	A	188	GLU	N-CA-C	-5.43	102.36	110.23
1	C	151	TRP	N-CA-C	-5.41	105.39	111.28
2	B	397	GLU	N-CA-C	5.39	122.27	110.80
1	C	204	LEU	N-CA-C	5.37	119.81	112.88
1	A	198	GLU	N-CA-C	-5.29	102.84	110.40
1	C	108	SER	N-CA-C	-5.27	105.54	111.28
1	A	756	SER	CA-C-N	5.25	125.71	120.04
1	A	756	SER	C-N-CA	5.25	125.71	120.04
2	D	653	LYS	N-CA-C	-5.25	105.56	111.28
2	B	309	ILE	CA-C-N	5.25	124.86	119.56
2	B	309	ILE	C-N-CA	5.25	124.86	119.56
2	D	407	ALA	N-CA-C	5.18	121.25	109.81
2	B	785	HIS	CA-C-N	5.18	128.18	120.31
2	B	785	HIS	C-N-CA	5.18	128.18	120.31
1	C	132	SER	N-CA-C	5.17	117.32	111.11
1	A	486	GLU	N-CA-C	5.16	115.05	108.24
2	D	37	VAL	CA-C-N	5.15	133.05	120.28
2	D	37	VAL	C-N-CA	5.15	133.05	120.28
2	D	106	ILE	N-CA-C	-5.14	105.38	110.62
1	A	137	PHE	N-CA-C	5.10	117.07	108.96
2	B	209	SER	N-CA-C	-5.10	104.60	111.54
1	A	255	SER	N-CA-C	5.09	117.43	110.55
1	C	744	PHE	N-CA-C	-5.09	105.10	111.92
2	B	765	GLN	N-CA-C	-5.07	105.84	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	220	SER	CA-C-N	5.06	124.56	119.19
2	B	220	SER	C-N-CA	5.06	124.56	119.19
2	D	669	VAL	CA-C-N	5.01	125.02	120.21
2	D	669	VAL	C-N-CA	5.01	125.02	120.21

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	542	LYS	Peptide
2	B	246	GLY	Peptide
2	B	409	PHE	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	2632	0	705	9	0
1	C	2508	0	664	10	0
2	B	2532	0	686	10	0
2	D	2509	0	689	12	0
All	All	10181	0	2744	41	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 (41) 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:C:660:PRO:O	1:C:664:ASN:N	2.15	0.79
2:D:50:GLU:O	2:D:53:ASP:N	2.20	0.74
1:A:660:PRO:O	1:A:664:ASN:N	2.22	0.73
1:A:464:LEU:O	1:A:468:MET:N	2.23	0.72
1:C:464:LEU:O	1:C:468:MET:N	2.20	0.72
1:A:483:GLY:HA2	1:A:497:ASN:O	1.93	0.68

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:47:ASP:O	2:D:51:LYS:N	2.29	0.65
2:D:330:LEU:O	2:D:334:LEU:N	2.31	0.64
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
1:C:265:GLY:HA3	1:C:381:TRP:C	2.23	0.62
2:D:406:GLU:N	2:D:410:VAL:O	2.33	0.62
2:D:493:GLU:O	2:D:499:ALA:N	2.31	0.60
2:D:526:GLY:O	2:D:716:ALA:N	2.36	0.58
1:C:102:LEU:O	1:C:105:THR:N	2.37	0.57
2:B:526:GLY:O	2:B:716:ALA:N	2.37	0.57
1:A:498:GLY:O	1:A:501:GLY:N	2.38	0.57
2:D:174:TYR:C	2:D:176:ASP:H	2.12	0.57
2:B:493:GLU:O	2:B:499:ALA:N	2.37	0.56
1:C:149:LEU:O	1:C:153:GLU:N	2.35	0.55
1:A:365:GLY:HA2	1:A:375:ASN:H	1.73	0.54
1:C:150:VAL:O	1:C:154:MET:N	2.35	0.53
2:D:303:LEU:O	2:D:307:ASN:N	2.36	0.53
1:C:365:GLY:HA2	1:C:375:ASN:H	1.75	0.52
2:B:303:LEU:O	2:B:307:ASN:N	2.34	0.52
1:A:30:GLY:HA2	1:A:63:THR:O	2.11	0.51
2:B:480:GLY:HA2	2:B:488:ASN:O	2.12	0.50
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	2.23	0.45
2:D:648:GLY:C	2:D:650:SER:H	2.25	0.44
1:A:490:ASN:O	1:A:491:SER:C	2.60	0.43
1:C:383:GLY:C	1:C:385:GLU:H	2.28	0.42
2:D:50:GLU:O	2:D:51:LYS:C	2.63	0.42
2:B:84:MET:O	2:B:85:SER:C	2.63	0.42
1:C:123:THR:C	1:C:125:MET:H	2.29	0.41
1:A:482:PHE:O	1:A:499:MET:N	2.53	0.41
2:B:100:GLN:O	2:B:124:GLY:HA3	2.20	0.41
1:C:265:GLY:HA3	1:C:381:TRP:O	2.21	0.40
1:A:710:ALA:O	1:A:715:LYS:N	2.54	0.40
2:B:761:LEU:O	2:B:762:ALA:C	2.64	0.40

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	652/822 (79%)	626 (96%)	22 (3%)	4 (1%)	21	59
1	C	618/822 (75%)	597 (97%)	20 (3%)	1 (0%)	43	78
2	B	625/825 (76%)	593 (95%)	28 (4%)	4 (1%)	21	59
2	D	619/825 (75%)	591 (96%)	27 (4%)	1 (0%)	43	78
All	All	2514/3294 (76%)	2407 (96%)	97 (4%)	10 (0%)	31	67

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	408	PRO
2	D	408	PRO
1	A	489	ASN
1	A	490	ASN
2	B	397	GLU
1	C	124	ARG
1	A	495	ALA
1	A	488	VAL
2	B	407	ALA
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 [i](#)

There are no chain breaks in this entry.

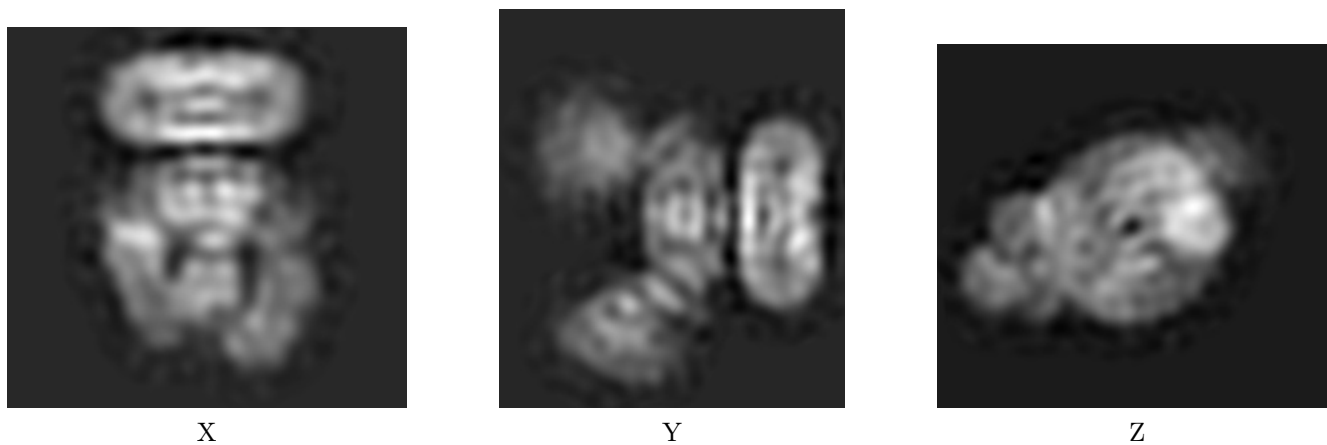
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8103. 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](#)

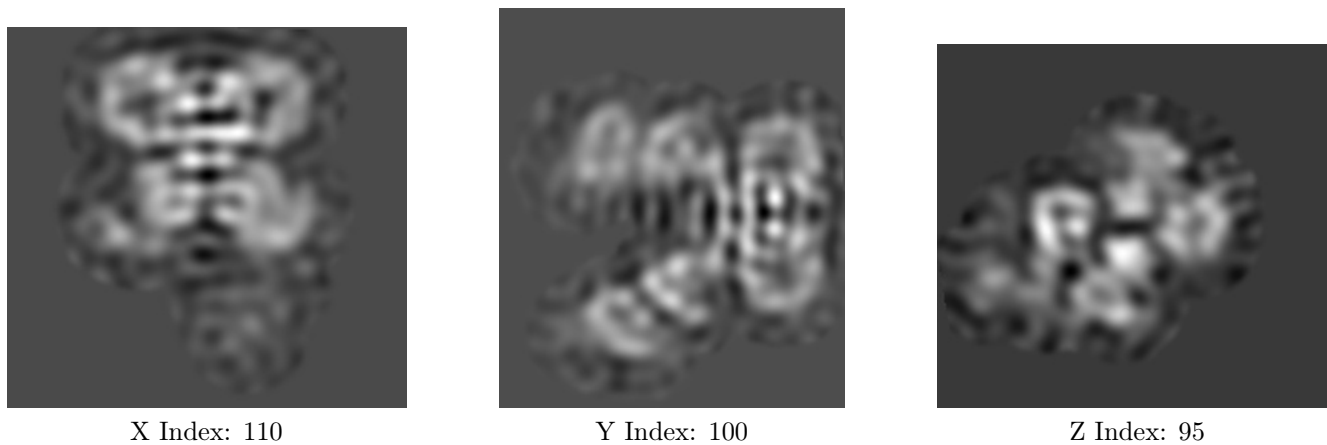
#### 6.1.1 Primary map



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](#)

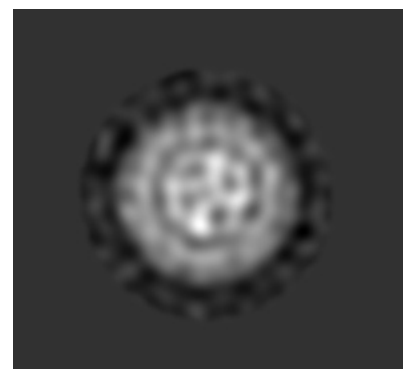
### 6.3.1 Primary map



X Index: 106



Y Index: 108



Z Index: 138

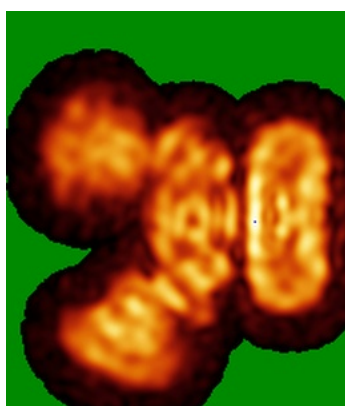
The images above show the largest variance slices of the map in three orthogonal directions.

## 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

This section was not generated.

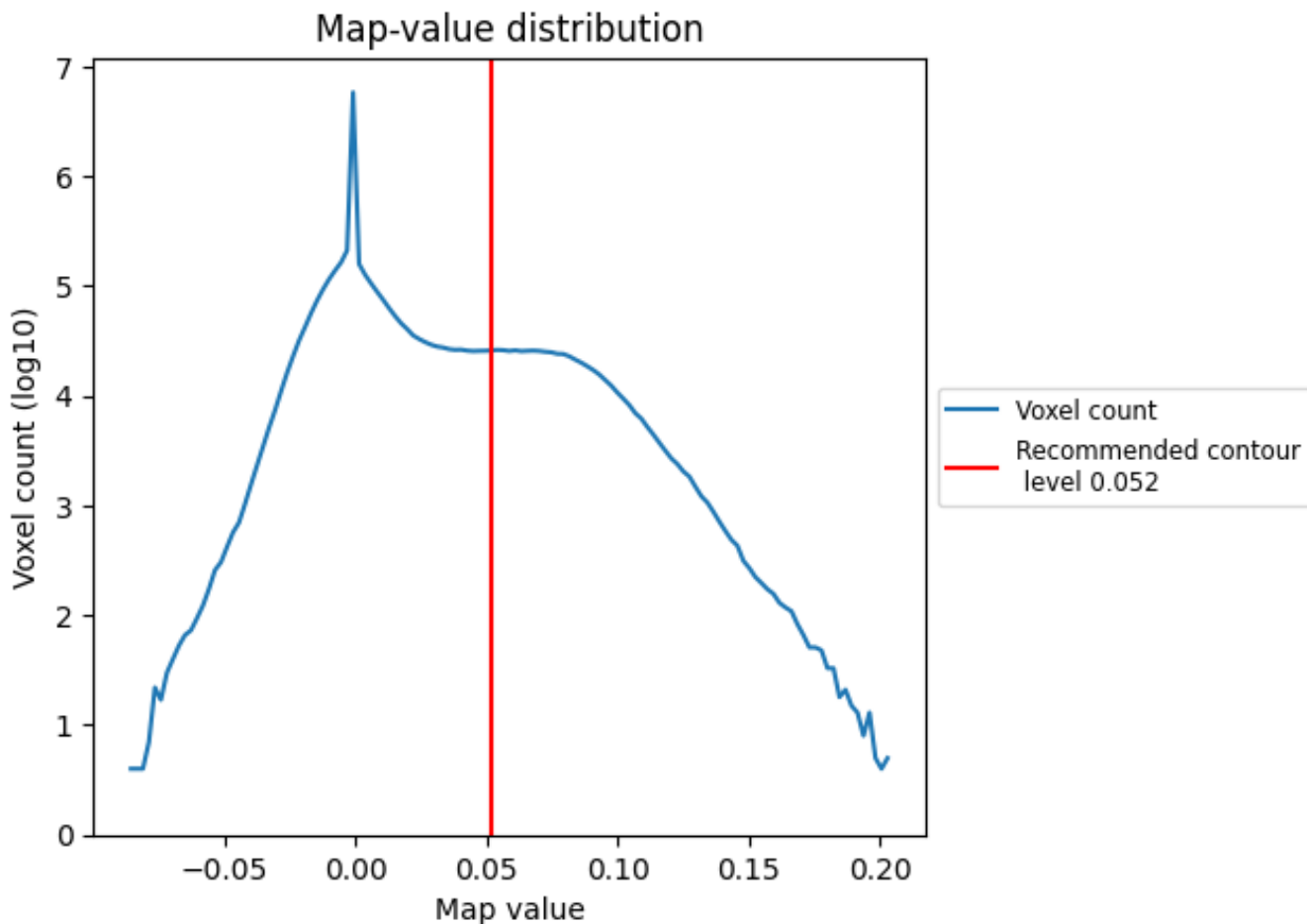
## 6.6 Mask visualisation

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

## 7 Map analysis [i](#)

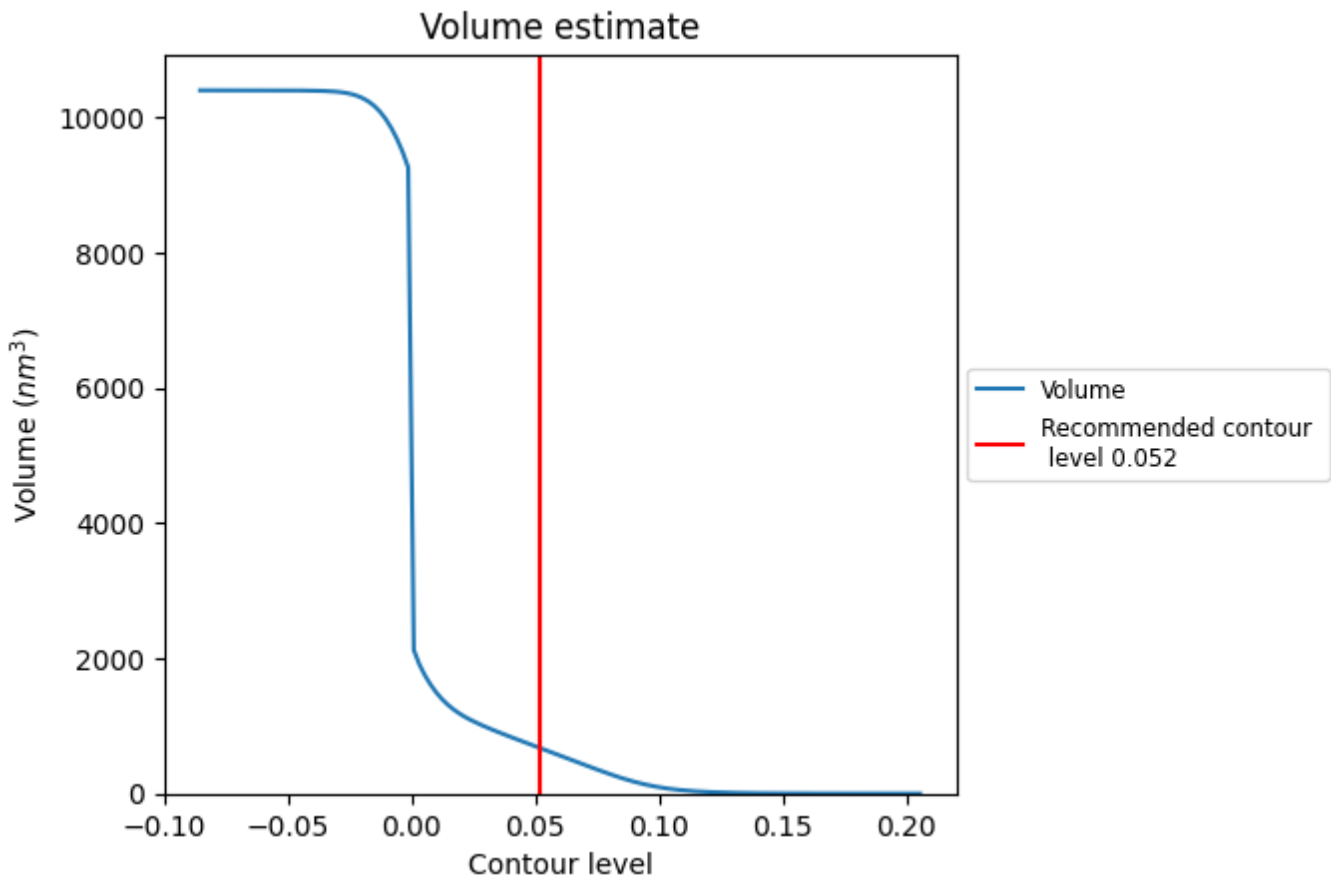
This section contains the results of statistical analysis of the map.

### 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 671 nm<sup>3</sup>; this corresponds to an approximate mass of 606 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.

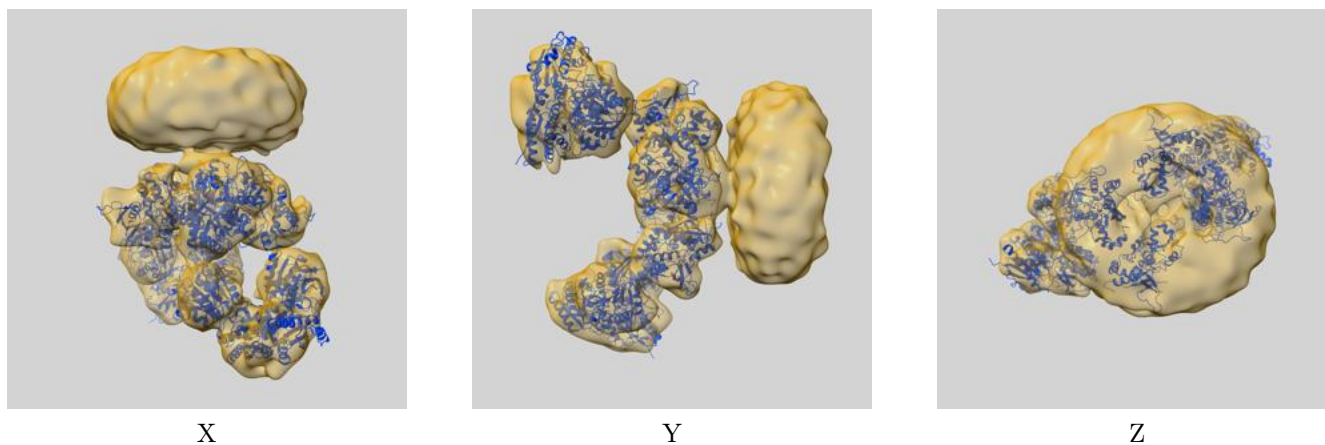
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

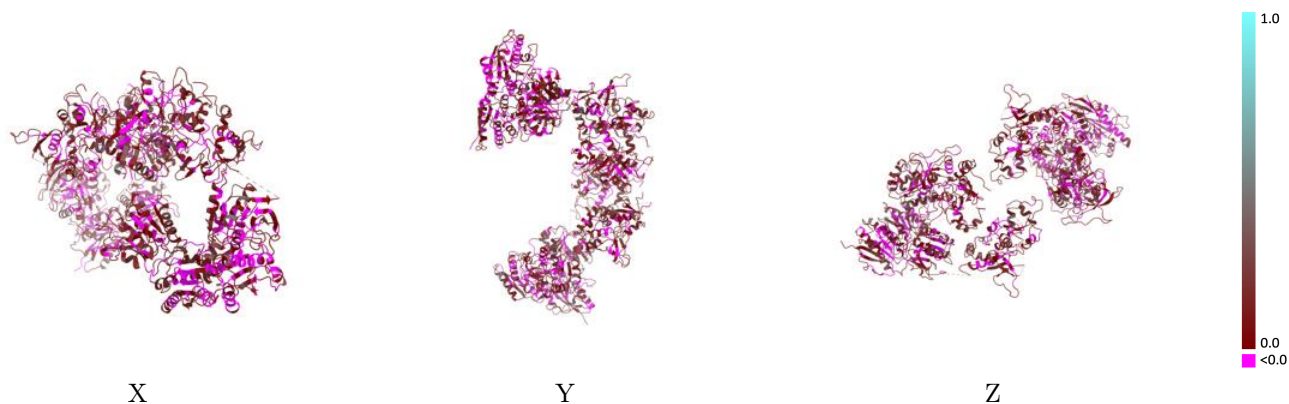
This section contains information regarding the fit between EMDB map EMD-8103 and PDB model 5IPS. Per-residue inclusion information can be found in section 3 on page 7.

### 9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.052 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](#)

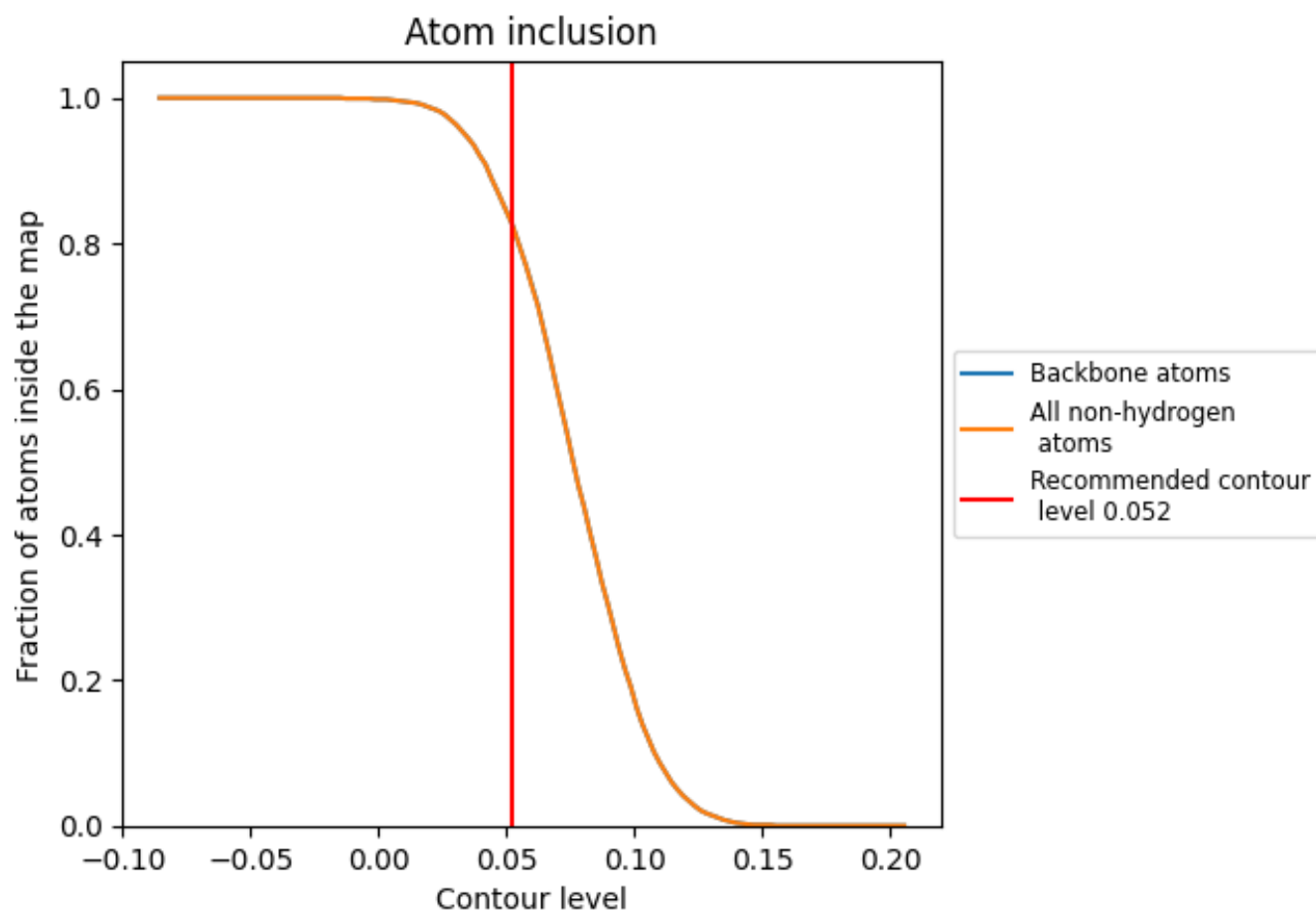


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.










## 9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 83% 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	 0.8280	 0.0890
A	 0.8970	 0.1060
B	 0.8240	 0.0870
C	 0.8310	 0.0890
D	 0.7560	 0.0720

