



Full wwPDB EM Validation Report ⓘ

Sep 19, 2026 – 07:29 am BST

PDB ID : 9SLM / pdb_00009slm
EMDB ID : EMD-55007
Title : GABA-A a4b3d receptor in complex with GABA and Nb24
Authors : Nestorow, S.; Miller, P.S.
Deposited on : 2025-09-04
Resolution : 3.27 Å(reported)
Based on initial model : 7QN7

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:

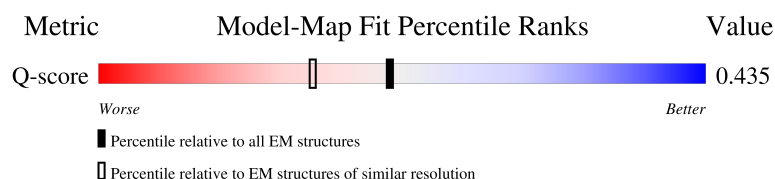
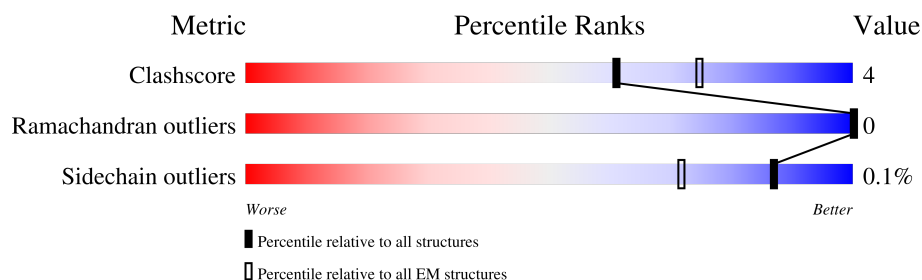
EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMD 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) : 2.50

1 Overall quality at a glance

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

The reported resolution of this entry is 3.27 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.



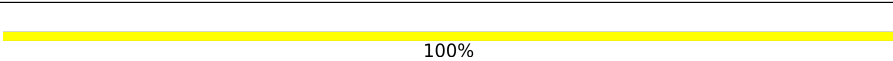
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14508 (2.77 - 3.77)

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	364	 81% 11% 8%
1	C	364	 85% 7% 8%
2	B	451	 67% 6% 26%
2	D	451	 63% 10% 27%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	E	430	
4	F	123	
4	G	123	
5	O	2	
6	P	3	

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 30946 atoms, of which 15325 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 Gamma-aminobutyric acid receptor subunit alpha-4, Gamma-aminobutyric acid type A receptor subunit alpha4.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	335	Total	C	H	N	O	S	0	0
			5416	1762	2714	437	483	20		
1	C	335	Total	C	H	N	O	S	0	0
			5418	1762	2716	437	483	20		

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q9D6F4
A	120	THR	MET	engineered mutation	UNP Q9D6F4
A	309	THR	-	linker	UNP Q9D6F4
A	310	ASN	-	linker	UNP Q9D6F4
A	311	SER	-	linker	UNP Q9D6F4
A	312	GLN	-	linker	UNP Q9D6F4
A	313	PRO	-	linker	UNP Q9D6F4
A	314	ALA	-	linker	UNP Q9D6F4
A	315	ARG	-	linker	UNP Q9D6F4
A	316	ALA	-	linker	UNP Q9D6F4
A	317	ALA	-	linker	UNP Q9D6F4
A	428	GLY	-	expression tag	UNP F6UBA8
A	429	THR	-	expression tag	UNP F6UBA8
A	430	LYS	-	expression tag	UNP F6UBA8
A	431	HIS	-	expression tag	UNP F6UBA8
A	432	HIS	-	expression tag	UNP F6UBA8
A	433	HIS	-	expression tag	UNP F6UBA8
A	434	HIS	-	expression tag	UNP F6UBA8
A	435	HIS	-	expression tag	UNP F6UBA8
A	436	HIS	-	expression tag	UNP F6UBA8
C	1	GLY	-	expression tag	UNP Q9D6F4
C	120	THR	MET	engineered mutation	UNP Q9D6F4
C	309	THR	-	linker	UNP Q9D6F4
C	310	ASN	-	linker	UNP Q9D6F4
C	311	SER	-	linker	UNP Q9D6F4

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	312	GLN	-	linker	UNP Q9D6F4
C	313	PRO	-	linker	UNP Q9D6F4
C	314	ALA	-	linker	UNP Q9D6F4
C	315	ARG	-	linker	UNP Q9D6F4
C	316	ALA	-	linker	UNP Q9D6F4
C	317	ALA	-	linker	UNP Q9D6F4
C	428	GLY	-	expression tag	UNP F6UBA8
C	429	THR	-	expression tag	UNP F6UBA8
C	430	LYS	-	expression tag	UNP F6UBA8
C	431	HIS	-	expression tag	UNP F6UBA8
C	432	HIS	-	expression tag	UNP F6UBA8
C	433	HIS	-	expression tag	UNP F6UBA8
C	434	HIS	-	expression tag	UNP F6UBA8
C	435	HIS	-	expression tag	UNP F6UBA8
C	436	HIS	-	expression tag	UNP F6UBA8

- Molecule 2 is a protein called Gamma-aminobutyric acid receptor subunit beta-3.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	332	Total	C	H	N	O	S	0	0
			5397	1780	2679	444	478	16		
2	D	331	Total	C	H	N	O	S	0	0
			5396	1778	2682	443	477	16		

There are 258 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLN	-	expression tag	UNP P28472
B	2	SER	-	expression tag	UNP P28472
B	3	VAL	-	expression tag	UNP P28472
B	4	ASN	-	expression tag	UNP P28472
B	5	ASP	-	expression tag	UNP P28472
B	6	PRO	-	expression tag	UNP P28472
B	302	VAL	VAL	linker	UNP P28472
B	303	ASN	ASN	linker	UNP P28472
B	304	TYR	TYR	linker	UNP P28472
B	305	ILE	ILE	linker	UNP P28472
B	306	PHE	PHE	linker	UNP P28472
B	307	PHE	PHE	linker	UNP P28472
B	307A	SER	-	linker	UNP P28472
B	307B	GLN	-	linker	UNP P28472
B	307C	PRO	-	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	307D	ALA	-	linker	UNP P28472
B	307E	GLY	-	linker	UNP P28472
B	307F	THR	-	linker	UNP P28472
B	307G	ALA	-	linker	UNP P28472
B	307H	ASP	-	linker	UNP P28472
B	307I	LEU	-	linker	UNP P28472
B	307J	GLU	-	linker	UNP P28472
B	307K	ASP	-	linker	UNP P28472
B	307L	ASN	-	linker	UNP P28472
B	307M	TRP	-	linker	UNP P28472
B	307N	GLU	-	linker	UNP P28472
B	307O	THR	-	linker	UNP P28472
B	307P	LEU	-	linker	UNP P28472
B	307Q	ASN	-	linker	UNP P28472
B	307R	ASP	-	linker	UNP P28472
B	307S	ASN	-	linker	UNP P28472
B	307T	LEU	-	linker	UNP P28472
B	307U	LYS	-	linker	UNP P28472
B	307V	VAL	-	linker	UNP P28472
B	307W	ILE	-	linker	UNP P28472
B	307X	GLU	-	linker	UNP P28472
B	307Y	LYS	-	linker	UNP P28472
B	307Z	ALA	-	linker	UNP P28472
B	308A	ASP	-	linker	UNP P28472
B	308B	ASN	-	linker	UNP P28472
B	308C	ALA	-	linker	UNP P28472
B	308D	ALA	-	linker	UNP P28472
B	308E	GLN	-	linker	UNP P28472
B	308F	VAL	-	linker	UNP P28472
B	308G	LYS	-	linker	UNP P28472
B	308H	ASP	-	linker	UNP P28472
B	308I	ALA	-	linker	UNP P28472
B	308J	LEU	-	linker	UNP P28472
B	308K	THR	-	linker	UNP P28472
B	308L	LYS	-	linker	UNP P28472
B	308M	MET	-	linker	UNP P28472
B	308N	ARG	-	linker	UNP P28472
B	308O	ALA	-	linker	UNP P28472
B	308P	ALA	-	linker	UNP P28472
B	308Q	ALA	-	linker	UNP P28472
B	308R	LEU	-	linker	UNP P28472
B	308S	ASP	-	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	308T	ALA	-	linker	UNP P28472
B	308U	GLN	-	linker	UNP P28472
B	308V	LYS	-	linker	UNP P28472
B	308W	ALA	-	linker	UNP P28472
B	308X	THR	-	linker	UNP P28472
B	308Y	PRO	-	linker	UNP P28472
B	308Z	PRO	-	linker	UNP P28472
B	309A	LYS	-	linker	UNP P28472
B	309B	LEU	-	linker	UNP P28472
B	309C	GLU	-	linker	UNP P28472
B	309D	ASP	-	linker	UNP P28472
B	309E	LYS	-	linker	UNP P28472
B	309F	SER	-	linker	UNP P28472
B	309G	PRO	-	linker	UNP P28472
B	309H	ASP	-	linker	UNP P28472
B	309I	SER	-	linker	UNP P28472
B	309J	PRO	-	linker	UNP P28472
B	309K	GLU	-	linker	UNP P28472
B	309L	MET	-	linker	UNP P28472
B	309M	LYS	-	linker	UNP P28472
B	309N	ASP	-	linker	UNP P28472
B	309O	PHE	-	linker	UNP P28472
B	309P	ARG	-	linker	UNP P28472
B	309Q	HIS	-	linker	UNP P28472
B	309R	GLY	-	linker	UNP P28472
B	309S	PHE	-	linker	UNP P28472
B	309T	ASP	-	linker	UNP P28472
B	309U	ILE	-	linker	UNP P28472
B	309V	LEU	-	linker	UNP P28472
B	309W	VAL	-	linker	UNP P28472
B	309X	GLY	-	linker	UNP P28472
B	309Y	GLN	-	linker	UNP P28472
B	309Z	ILE	-	linker	UNP P28472
B	310A	ASP	-	linker	UNP P28472
B	310B	ASP	-	linker	UNP P28472
B	310C	ALA	-	linker	UNP P28472
B	310D	LEU	-	linker	UNP P28472
B	310E	LYS	-	linker	UNP P28472
B	310F	LEU	-	linker	UNP P28472
B	310G	ALA	-	linker	UNP P28472
B	310H	ASN	-	linker	UNP P28472
B	310I	GLU	-	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	310J	GLY	-	linker	UNP P28472
B	310K	LYS	-	linker	UNP P28472
B	310L	VAL	-	linker	UNP P28472
B	310M	LYS	-	linker	UNP P28472
B	310N	GLU	-	linker	UNP P28472
B	310O	ALA	-	linker	UNP P28472
B	310P	GLN	-	linker	UNP P28472
B	310Q	ALA	-	linker	UNP P28472
B	310R	ALA	-	linker	UNP P28472
B	310S	ALA	-	linker	UNP P28472
B	310T	GLU	-	linker	UNP P28472
B	310U	GLN	-	linker	UNP P28472
B	310V	LEU	-	linker	UNP P28472
B	310W	LYS	-	linker	UNP P28472
B	310X	THR	-	linker	UNP P28472
B	310Y	THR	-	linker	UNP P28472
B	310Z	ARG	-	linker	UNP P28472
B	311A	ASN	-	linker	UNP P28472
B	311B	ALA	-	linker	UNP P28472
B	311C	TYR	-	linker	UNP P28472
B	311D	ILE	-	linker	UNP P28472
B	311E	GLN	-	linker	UNP P28472
B	311F	LYS	-	linker	UNP P28472
B	311G	TYR	-	linker	UNP P28472
B	311H	LEU	-	linker	UNP P28472
B	417	THR	-	linker	UNP P28472
B	418	GLY	-	linker	UNP P28472
B	419	ARG	-	linker	UNP P28472
B	420	ALA	-	linker	UNP P28472
B	421	ALA	-	linker	UNP P28472
D	1	GLN	-	expression tag	UNP P28472
D	2	SER	-	expression tag	UNP P28472
D	3	VAL	-	expression tag	UNP P28472
D	4	ASN	-	expression tag	UNP P28472
D	5	ASP	-	expression tag	UNP P28472
D	6	PRO	-	expression tag	UNP P28472
D	302	VAL	VAL	linker	UNP P28472
D	303	ASN	ASN	linker	UNP P28472
D	304	TYR	TYR	linker	UNP P28472
D	305	ILE	ILE	linker	UNP P28472
D	306	PHE	PHE	linker	UNP P28472
D	307	PHE	PHE	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	307A	SER	-	linker	UNP P28472
D	307B	GLN	-	linker	UNP P28472
D	307C	PRO	-	linker	UNP P28472
D	307D	ALA	-	linker	UNP P28472
D	307E	GLY	-	linker	UNP P28472
D	307F	THR	-	linker	UNP P28472
D	307G	ALA	-	linker	UNP P28472
D	307H	ASP	-	linker	UNP P28472
D	307I	LEU	-	linker	UNP P28472
D	307J	GLU	-	linker	UNP P28472
D	307K	ASP	-	linker	UNP P28472
D	307L	ASN	-	linker	UNP P28472
D	307M	TRP	-	linker	UNP P28472
D	307N	GLU	-	linker	UNP P28472
D	307O	THR	-	linker	UNP P28472
D	307P	LEU	-	linker	UNP P28472
D	307Q	ASN	-	linker	UNP P28472
D	307R	ASP	-	linker	UNP P28472
D	307S	ASN	-	linker	UNP P28472
D	307T	LEU	-	linker	UNP P28472
D	307U	LYS	-	linker	UNP P28472
D	307V	VAL	-	linker	UNP P28472
D	307W	ILE	-	linker	UNP P28472
D	307X	GLU	-	linker	UNP P28472
D	307Y	LYS	-	linker	UNP P28472
D	307Z	ALA	-	linker	UNP P28472
D	308A	ASP	-	linker	UNP P28472
D	308B	ASN	-	linker	UNP P28472
D	308C	ALA	-	linker	UNP P28472
D	308D	ALA	-	linker	UNP P28472
D	308E	GLN	-	linker	UNP P28472
D	308F	VAL	-	linker	UNP P28472
D	308G	LYS	-	linker	UNP P28472
D	308H	ASP	-	linker	UNP P28472
D	308I	ALA	-	linker	UNP P28472
D	308J	LEU	-	linker	UNP P28472
D	308K	THR	-	linker	UNP P28472
D	308L	LYS	-	linker	UNP P28472
D	308M	MET	-	linker	UNP P28472
D	308N	ARG	-	linker	UNP P28472
D	308O	ALA	-	linker	UNP P28472
D	308P	ALA	-	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	308Q	ALA	-	linker	UNP P28472
D	308R	LEU	-	linker	UNP P28472
D	308S	ASP	-	linker	UNP P28472
D	308T	ALA	-	linker	UNP P28472
D	308U	GLN	-	linker	UNP P28472
D	308V	LYS	-	linker	UNP P28472
D	308W	ALA	-	linker	UNP P28472
D	308X	THR	-	linker	UNP P28472
D	308Y	PRO	-	linker	UNP P28472
D	308Z	PRO	-	linker	UNP P28472
D	309A	LYS	-	linker	UNP P28472
D	309B	LEU	-	linker	UNP P28472
D	309C	GLU	-	linker	UNP P28472
D	309D	ASP	-	linker	UNP P28472
D	309E	LYS	-	linker	UNP P28472
D	309F	SER	-	linker	UNP P28472
D	309G	PRO	-	linker	UNP P28472
D	309H	ASP	-	linker	UNP P28472
D	309I	SER	-	linker	UNP P28472
D	309J	PRO	-	linker	UNP P28472
D	309K	GLU	-	linker	UNP P28472
D	309L	MET	-	linker	UNP P28472
D	309M	LYS	-	linker	UNP P28472
D	309N	ASP	-	linker	UNP P28472
D	309O	PHE	-	linker	UNP P28472
D	309P	ARG	-	linker	UNP P28472
D	309Q	HIS	-	linker	UNP P28472
D	309R	GLY	-	linker	UNP P28472
D	309S	PHE	-	linker	UNP P28472
D	309T	ASP	-	linker	UNP P28472
D	309U	ILE	-	linker	UNP P28472
D	309V	LEU	-	linker	UNP P28472
D	309W	VAL	-	linker	UNP P28472
D	309X	GLY	-	linker	UNP P28472
D	309Y	GLN	-	linker	UNP P28472
D	309Z	ILE	-	linker	UNP P28472
D	310A	ASP	-	linker	UNP P28472
D	310B	ASP	-	linker	UNP P28472
D	310C	ALA	-	linker	UNP P28472
D	310D	LEU	-	linker	UNP P28472
D	310E	LYS	-	linker	UNP P28472
D	310F	LEU	-	linker	UNP P28472

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	310G	ALA	-	linker	UNP P28472
D	310H	ASN	-	linker	UNP P28472
D	310I	GLU	-	linker	UNP P28472
D	310J	GLY	-	linker	UNP P28472
D	310K	LYS	-	linker	UNP P28472
D	310L	VAL	-	linker	UNP P28472
D	310M	LYS	-	linker	UNP P28472
D	310N	GLU	-	linker	UNP P28472
D	310O	ALA	-	linker	UNP P28472
D	310P	GLN	-	linker	UNP P28472
D	310Q	ALA	-	linker	UNP P28472
D	310R	ALA	-	linker	UNP P28472
D	310S	ALA	-	linker	UNP P28472
D	310T	GLU	-	linker	UNP P28472
D	310U	GLN	-	linker	UNP P28472
D	310V	LEU	-	linker	UNP P28472
D	310W	LYS	-	linker	UNP P28472
D	310X	THR	-	linker	UNP P28472
D	310Y	THR	-	linker	UNP P28472
D	310Z	ARG	-	linker	UNP P28472
D	311A	ASN	-	linker	UNP P28472
D	311B	ALA	-	linker	UNP P28472
D	311C	TYR	-	linker	UNP P28472
D	311D	ILE	-	linker	UNP P28472
D	311E	GLN	-	linker	UNP P28472
D	311F	LYS	-	linker	UNP P28472
D	311G	TYR	-	linker	UNP P28472
D	311H	LEU	-	linker	UNP P28472
D	417	THR	-	linker	UNP P28472
D	418	GLY	-	linker	UNP P28472
D	419	ARG	-	linker	UNP P28472
D	420	ALA	-	linker	UNP P28472
D	421	ALA	-	linker	UNP P28472

- Molecule 3 is a protein called Gamma-aminobutyric acid receptor subunit delta.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	E	328	Total	C	H	N	O	S	0	0
			5183	1698	2564	434	474	13		

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-83	MET	-	initiating methionine	UNP O14764
E	-82	GLY	-	expression tag	UNP O14764
E	-81	ILE	-	expression tag	UNP O14764
E	-80	LEU	-	expression tag	UNP O14764
E	-79	PRO	-	expression tag	UNP O14764
E	-78	SER	-	expression tag	UNP O14764
E	-77	PRO	-	expression tag	UNP O14764
E	-76	GLY	-	expression tag	UNP O14764
E	-75	MET	-	expression tag	UNP O14764
E	-74	PRO	-	expression tag	UNP O14764
E	-73	ALA	-	expression tag	UNP O14764
E	-72	LEU	-	expression tag	UNP O14764
E	-71	LEU	-	expression tag	UNP O14764
E	-70	SER	-	expression tag	UNP O14764
E	-69	LEU	-	expression tag	UNP O14764
E	-68	VAL	-	expression tag	UNP O14764
E	-67	SER	-	expression tag	UNP O14764
E	-66	LEU	-	expression tag	UNP O14764
E	-65	LEU	-	expression tag	UNP O14764
E	-64	SER	-	expression tag	UNP O14764
E	-63	VAL	-	expression tag	UNP O14764
E	-62	LEU	-	expression tag	UNP O14764
E	-61	LEU	-	expression tag	UNP O14764
E	-60	MET	-	expression tag	UNP O14764
E	-59	GLY	-	expression tag	UNP O14764
E	-58	CYS	-	expression tag	UNP O14764
E	-57	VAL	-	expression tag	UNP O14764
E	-56	ALA	-	expression tag	UNP O14764
E	-55	GLU	-	expression tag	UNP O14764
E	-54	MET	-	expression tag	UNP O14764
E	-53	ASP	-	expression tag	UNP O14764
E	-52	GLU	-	expression tag	UNP O14764
E	-51	LYS	-	expression tag	UNP O14764
E	-50	THR	-	expression tag	UNP O14764
E	-49	THR	-	expression tag	UNP O14764
E	-48	GLY	-	expression tag	UNP O14764
E	-47	TRP	-	expression tag	UNP O14764
E	-46	ARG	-	expression tag	UNP O14764
E	-45	GLY	-	expression tag	UNP O14764
E	-44	GLY	-	expression tag	UNP O14764
E	-43	HIS	-	expression tag	UNP O14764
E	-42	VAL	-	expression tag	UNP O14764
E	-41	VAL	-	expression tag	UNP O14764

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	-40	GLU	-	expression tag	UNP O14764
E	-39	GLY	-	expression tag	UNP O14764
E	-38	LEU	-	expression tag	UNP O14764
E	-37	ALA	-	expression tag	UNP O14764
E	-36	GLY	-	expression tag	UNP O14764
E	-35	GLU	-	expression tag	UNP O14764
E	-34	LEU	-	expression tag	UNP O14764
E	-33	GLU	-	expression tag	UNP O14764
E	-32	GLN	-	expression tag	UNP O14764
E	-31	LEU	-	expression tag	UNP O14764
E	-30	ARG	-	expression tag	UNP O14764
E	-29	ALA	-	expression tag	UNP O14764
E	-28	ARG	-	expression tag	UNP O14764
E	-27	LEU	-	expression tag	UNP O14764
E	-26	GLU	-	expression tag	UNP O14764
E	-25	HIS	-	expression tag	UNP O14764
E	-24	HIS	-	expression tag	UNP O14764
E	-23	PRO	-	expression tag	UNP O14764
E	-22	GLN	-	expression tag	UNP O14764
E	-21	GLY	-	expression tag	UNP O14764
E	-20	GLN	-	expression tag	UNP O14764
E	-19	ARG	-	expression tag	UNP O14764
E	-18	GLU	-	expression tag	UNP O14764
E	-17	PRO	-	expression tag	UNP O14764
E	-16	ASP	-	expression tag	UNP O14764
E	-15	TYR	-	expression tag	UNP O14764
E	-14	ASP	-	expression tag	UNP O14764
E	-13	ILE	-	expression tag	UNP O14764
E	-12	PRO	-	expression tag	UNP O14764
E	-11	THR	-	expression tag	UNP O14764
E	-10	THR	-	expression tag	UNP O14764
E	-9	GLU	-	expression tag	UNP O14764
E	-8	ASN	-	expression tag	UNP O14764
E	-7	LEU	-	expression tag	UNP O14764
E	-6	TYR	-	expression tag	UNP O14764
E	-5	PHE	-	expression tag	UNP O14764
E	-4	GLN	-	expression tag	UNP O14764
E	-3	GLY	-	expression tag	UNP O14764
E	-2	THR	-	expression tag	UNP O14764
E	-1	GLY	-	expression tag	UNP O14764
E	219	LEU	-	linker	UNP O14764
E	220	ARG	-	linker	UNP O14764

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	221	ARG	-	linker	UNP O14764
E	222	ASN	-	linker	UNP O14764
E	223	ARG	-	linker	UNP O14764
E	224	GLY	-	linker	UNP O14764
E	225	VAL	-	linker	UNP O14764
E	226	TYR	-	linker	UNP O14764
E	227	ILE	-	linker	UNP O14764
E	228	ILE	-	linker	UNP O14764
E	229	GLN	-	linker	UNP O14764
E	230	SER	-	linker	UNP O14764
E	231	TYR	-	linker	UNP O14764
E	232	MET	-	linker	UNP O14764
E	233	PRO	-	linker	UNP O14764
E	234	SER	-	linker	UNP O14764
E	235	VAL	-	linker	UNP O14764
E	236	LEU	-	linker	UNP O14764
E	237	LEU	-	linker	UNP O14764
E	238	VAL	-	linker	UNP O14764
E	239	ALA	-	linker	UNP O14764
E	240	MET	-	linker	UNP O14764
E	241	SER	-	linker	UNP O14764
E	242	TRP	-	linker	UNP O14764
E	243	VAL	-	linker	UNP O14764
E	244	SER	-	linker	UNP O14764
E	245	PHE	-	linker	UNP O14764
E	246	TRP	-	linker	UNP O14764
E	247	ILE	-	linker	UNP O14764
E	248	SER	-	linker	UNP O14764
E	249	GLN	-	linker	UNP O14764
E	250	ALA	-	linker	UNP O14764
E	251	ALA	-	linker	UNP O14764
E	252	VAL	-	linker	UNP O14764
E	253	PRO	-	linker	UNP O14764
E	254	ALA	-	linker	UNP O14764
E	255	ARG	-	linker	UNP O14764
E	256	VAL	-	linker	UNP O14764
E	257	SER	-	linker	UNP O14764
E	258	LEU	-	linker	UNP O14764
E	259	GLY	-	linker	UNP O14764
E	260	ILE	-	linker	UNP O14764
E	261	THR	-	linker	UNP O14764
E	262	THR	-	linker	UNP O14764

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	263	VAL	-	linker	UNP O14764
E	264	LEU	-	linker	UNP O14764
E	265	THR	-	linker	UNP O14764
E	266	MET	-	linker	UNP O14764
E	267	THR	-	linker	UNP O14764
E	268	THR	-	linker	UNP O14764
E	269	LEU	-	linker	UNP O14764
E	270	MET	-	linker	UNP O14764
E	271	VAL	-	linker	UNP O14764
E	272	SER	-	linker	UNP O14764
E	273	ALA	-	linker	UNP O14764
E	274	ARG	-	linker	UNP O14764
E	275	SER	-	linker	UNP O14764
E	276	SER	-	linker	UNP O14764
E	277	LEU	-	linker	UNP O14764
E	278	PRO	-	linker	UNP O14764
E	279	ARG	-	linker	UNP O14764
E	280	ALA	-	linker	UNP O14764
E	281	SER	-	linker	UNP O14764
E	282	ALA	-	linker	UNP O14764
E	283	ILE	-	linker	UNP O14764
E	284	LYS	-	linker	UNP O14764
E	285	ALA	-	linker	UNP O14764
E	286	LEU	-	linker	UNP O14764
E	287	ASP	-	linker	UNP O14764
E	288	VAL	-	linker	UNP O14764
E	289	TYR	-	linker	UNP O14764
E	290	PHE	-	linker	UNP O14764
E	291	TRP	-	linker	UNP O14764
E	292	ILE	-	linker	UNP O14764
E	293	CYS	-	linker	UNP O14764
E	294	TYR	-	linker	UNP O14764
E	295	VAL	-	linker	UNP O14764
E	296	PHE	-	linker	UNP O14764
E	297	VAL	-	linker	UNP O14764
E	298	PHE	-	linker	UNP O14764
E	299	ALA	-	linker	UNP O14764
E	300	ALA	-	linker	UNP O14764
E	301	LEU	-	linker	UNP O14764
E	302	VAL	-	linker	UNP O14764
E	303	GLU	-	linker	UNP O14764
E	304	TYR	-	linker	UNP O14764

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	305	ALA	-	linker	UNP O14764
E	306	PHE	-	linker	UNP O14764
E	307	ALA	-	linker	UNP O14764
E	308	HIS	-	linker	UNP O14764
E	309	PHE	-	linker	UNP O14764
E	310	ASN	-	linker	UNP O14764
E	311	ALA	-	linker	UNP O14764
E	312	ASP	-	linker	UNP O14764
E	313	SER	-	linker	UNP O14764
E	314	GLN	-	linker	UNP O14764
E	315	PRO	-	linker	UNP O14764
E	316	ALA	-	linker	UNP O14764
E	317	ARG	-	linker	UNP O14764
E	318	ALA	-	linker	UNP O14764
E	319	ALA	-	linker	UNP O14764

- Molecule 4 is a protein called Nanobody Nb24.

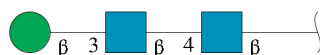
Mol	Chain	Residues	Atoms						AltConf	Trace
4	F	123	Total	C	H	N	O	S	0	0
			1841	583	896	171	187	4		
4	G	123	Total	C	H	N	O	S	0	0
			1841	583	896	171	187	4		

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



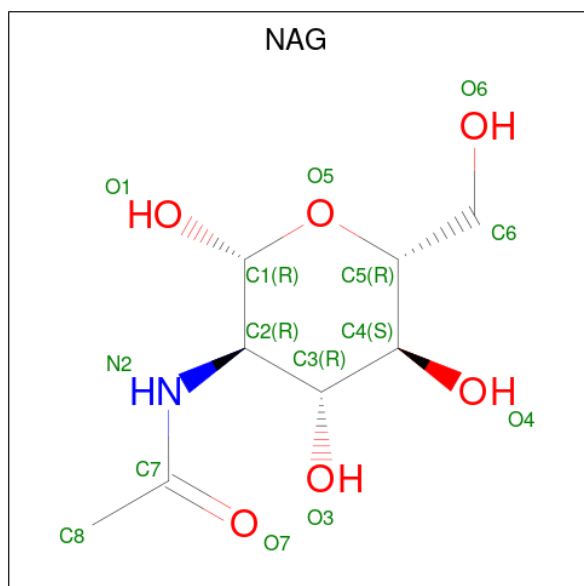
Mol	Chain	Residues	Atoms					AltConf	Trace
5	O	2	Total	C	H	N	O	0	0
			45	14	20	1	10		

- Molecule 6 is an oligosaccharide called beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



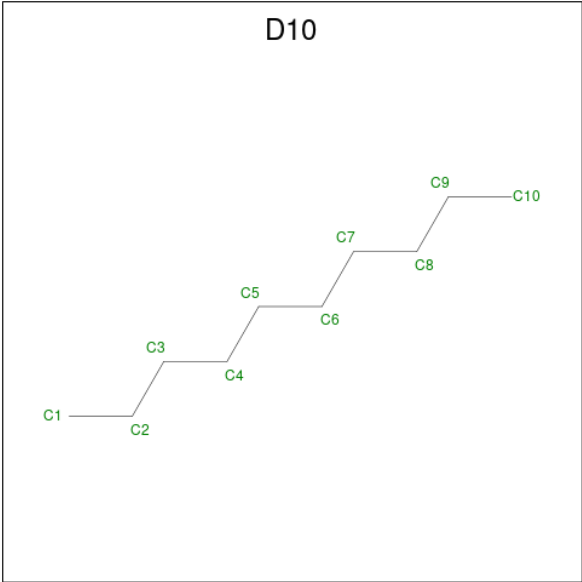
Mol	Chain	Residues	Atoms					AltConf	Trace
6	P	3	Total	C	H	N	O	0	0
			71	22	32	2	15		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



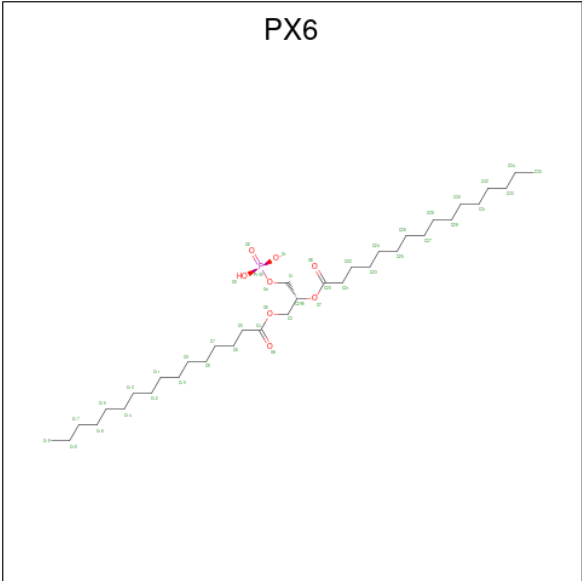
Mol	Chain	Residues	Atoms					AltConf
7	A	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	A	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	B	1	Total	C	H	N	O	0
			26	8	12	1	5	
7	B	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	B	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	C	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	D	1	Total	C	H	N	O	0
			28	8	14	1	5	
7	E	1	Total	C	H	N	O	0
			28	8	14	1	5	

- Molecule 8 is DECANE (CCD ID: D10) (formula: $C_{10}H_{22}$).



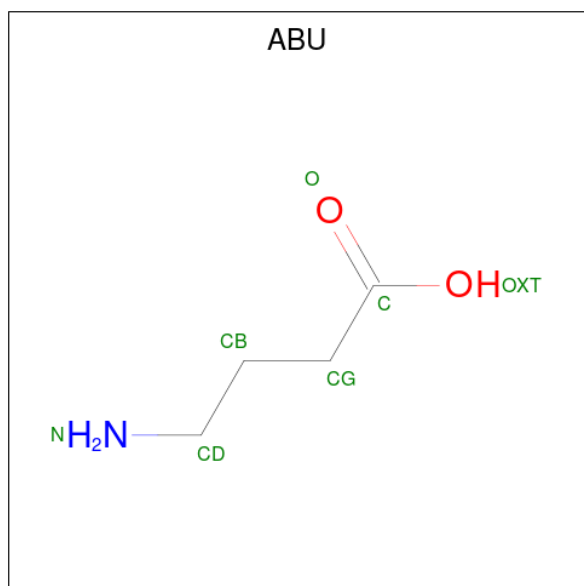
Mol	Chain	Residues	Atoms	AltConf
8	A	1	Total C 10 10	0
8	A	1	Total C 10 10	0
8	B	1	Total C 10 10	0
8	D	1	Total C 10 10	0

- Molecule 9 is 1,2-DIPALMITOYL-SN-GLYCERO-3-PHOSPHATE (CCD ID: PX6) (formula: C₃₅H₆₈O₈P).



Mol	Chain	Residues	Atoms				AltConf
9	A	1	Total	C	O	P	0
			44	35	8	1	

- Molecule 10 is GAMMA-AMINO-BUTANOIC ACID (CCD ID: ABU) (formula: $C_4H_9NO_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
10	B	1	Total	C	H	N	O	0
			15	4	8	1	2	
10	D	1	Total	C	H	N	O	0
			15	4	8	1	2	

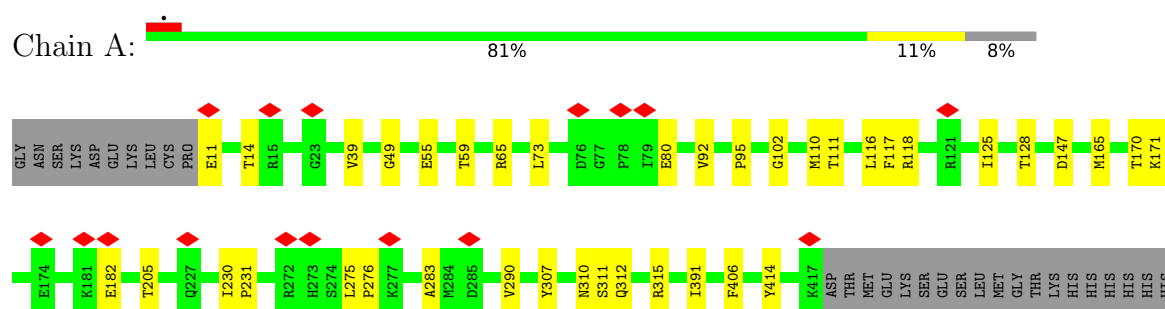
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		AltConf
11	A	1	Total	O	0
			1	1	
11	C	1	Total	O	0
			1	1	

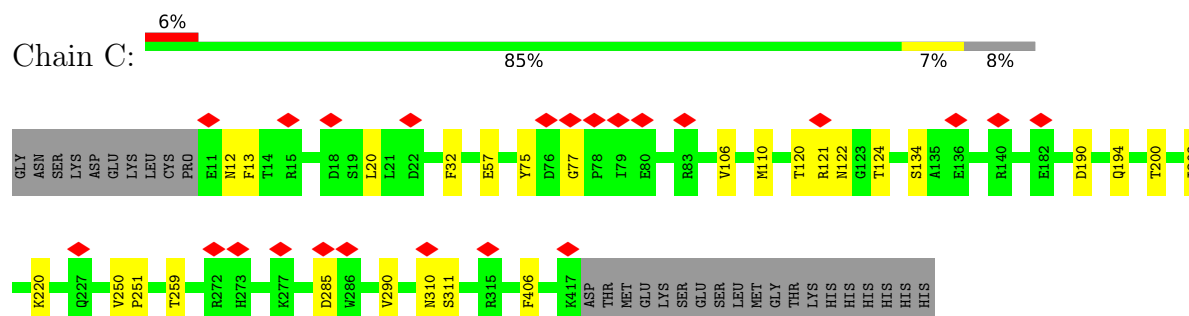
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

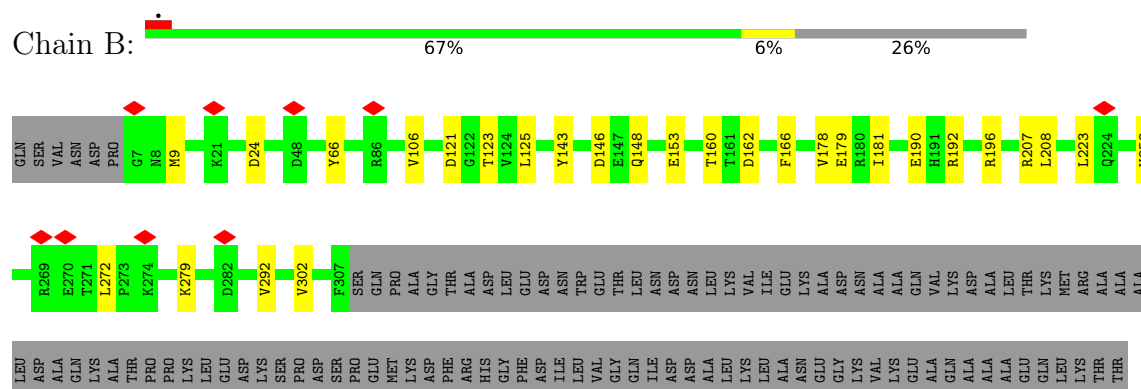
- Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-4, Gamma-aminobutyric acid type A receptor subunit alpha4

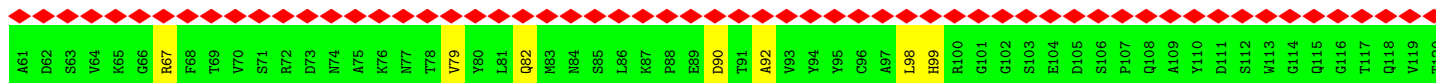


- Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-4, Gamma-aminobutyric acid type A receptor subunit alpha4



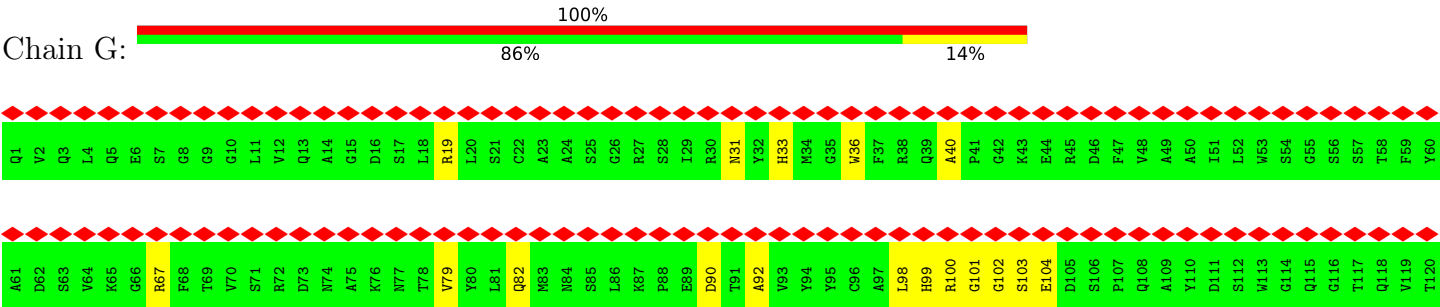
- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3







• Molecule 4: Nanobody Nb24



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 6: beta-D-mannopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	47269	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49.95	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	130000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.225	Depositor
Minimum map value	-0.134	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.0559	Depositor
Map size (Å)	247.68, 247.68, 247.68	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.645, 0.645, 0.645	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: D10, BMA, PX6, ABU, NAG

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	0.21	0/2775	0.34	0/3768
1	C	0.20	0/2775	0.33	0/3768
2	B	0.23	0/2790	0.34	0/3795
2	D	0.29	0/2786	0.46	0/3790
3	E	0.20	0/2692	0.32	0/3672
4	F	0.11	0/965	0.24	0/1303
4	G	0.11	0/965	0.25	0/1303
All	All	0.22	0/15748	0.35	0/21399

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2702	2714	2710	29	0
1	C	2702	2716	2711	21	0
2	B	2718	2679	2711	19	0
2	D	2714	2682	2709	41	0
3	E	2619	2564	2563	31	0
4	F	945	896	894	5	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	945	896	894	16	0
5	O	25	20	22	5	0
6	P	39	32	34	0	0
7	A	28	28	26	0	0
7	B	42	40	39	4	0
7	C	14	14	13	2	0
7	D	14	14	13	0	0
7	E	14	14	13	0	0
8	A	20	0	44	1	0
8	B	10	0	22	0	0
8	D	10	0	22	0	0
9	A	44	0	68	3	0
10	B	7	8	0	0	0
10	D	7	8	0	0	0
11	A	1	0	0	0	0
11	C	1	0	0	0	0
All	All	15621	15325	15508	138	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.

All (138) 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)
7:B:502:NAG:O3	5:O:1:NAG:C1	1.70	1.39
7:B:502:NAG:C3	5:O:1:NAG:C1	2.49	0.90
1:A:315:ARG:NE	9:A:504:PX6:O4	2.07	0.87
1:A:95:PRO:HB3	1:A:165:MET:HE1	1.62	0.82
2:D:253:LEU:HD23	3:E:260:ILE:HG21	1.63	0.80
2:D:190:GLU:OE2	2:D:213:ARG:NH1	2.21	0.73
1:C:77:GLY:O	1:C:121:ARG:NH2	2.27	0.68
7:B:502:NAG:HO3	5:O:1:NAG:C1	2.07	0.67
1:A:391:ILE:HD11	9:A:504:PX6:H3	1.76	0.67
1:C:190:ASP:OD1	1:C:220:LYS:NZ	2.28	0.66
3:E:197:PHE:CE1	3:E:215:LEU:HD13	2.31	0.66
2:D:70:LYS:O	4:G:103:SER:HA	1.95	0.65
2:D:43:ASP:OD1	2:D:176:THR:OG1	2.15	0.63
1:A:275:LEU:HD12	1:A:276:PRO:HD2	1.80	0.63
2:B:66:TYR:CE2	2:B:125:LEU:HD13	2.33	0.62
1:C:120:THR:OG1	1:C:124:THR:OG1	2.15	0.62
1:C:200:THR:HG22	1:C:209:ILE:HD13	1.79	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:253:LEU:HD21	3:E:256:VAL:HG13	1.81	0.61
4:F:19:ARG:NE	4:F:82:GLN:OE1	2.34	0.60
4:G:19:ARG:NE	4:G:82:GLN:OE1	2.34	0.60
1:A:92:VAL:HG21	1:A:117:PHE:CZ	2.38	0.59
2:B:223:LEU:HD21	1:C:285:ASP:HB3	1.84	0.59
3:E:159:GLU:OE2	3:E:203:ASN:ND2	2.36	0.58
1:A:80:GLU:N	1:A:80:GLU:OE1	2.33	0.58
3:E:78:TYR:OH	3:E:83:GLU:O	2.21	0.58
2:D:120:PRO:O	4:G:33:HIS:HE1	1.87	0.57
2:D:153:GLU:OE2	2:D:207:ARG:NH1	2.37	0.57
1:A:310:ASN:OD1	1:A:311:SER:N	2.37	0.57
2:B:146:ASP:OD2	2:B:148:GLN:NE2	2.38	0.57
1:A:116:LEU:HD12	1:A:117:PHE:H	1.69	0.57
1:C:110:MET:HE1	2:D:106:VAL:HG23	1.87	0.57
2:D:256:THR:HG21	3:E:260:ILE:HD12	1.87	0.56
3:E:174:GLU:N	3:E:174:GLU:OE1	2.38	0.56
3:E:92:VAL:HG22	3:E:120:ILE:CD1	2.37	0.55
7:B:502:NAG:H3	5:O:1:NAG:C1	2.37	0.55
3:E:92:VAL:HG22	3:E:120:ILE:HD12	1.87	0.55
3:E:150:ASP:OD2	3:E:152:GLN:NE2	2.40	0.55
4:G:67:ARG:NH2	4:G:90:ASP:OD2	2.40	0.55
1:C:310:ASN:OD1	1:C:311:SER:N	2.40	0.55
2:B:66:TYR:CZ	2:B:125:LEU:HD13	2.42	0.54
1:C:12:ASN:OD1	1:C:13:PHE:N	2.40	0.54
4:F:67:ARG:NH2	4:F:90:ASP:OD2	2.40	0.54
3:E:274:ARG:NH1	3:E:287:ASP:OD2	2.41	0.54
2:B:160:THR:OG1	2:B:162:ASP:OD1	2.26	0.53
2:D:68:ARG:NH2	4:G:104:GLU:OE1	2.41	0.53
2:D:88:ALA:HB2	2:D:116:ILE:HD11	1.88	0.53
1:A:311:SER:OG	1:A:312:GLN:OE1	2.17	0.53
2:D:305:ILE:HD12	2:D:420:ALA:CB	2.38	0.53
1:A:311:SER:HG	1:A:312:GLN:CD	2.13	0.53
2:D:233:THR:O	2:D:236:SER:OG	2.25	0.53
2:D:19:LEU:HD21	2:D:74:TYR:HB3	1.91	0.52
1:C:250:VAL:HB	1:C:251:PRO:HD3	1.92	0.52
2:D:264:ILE:HD13	3:E:271:VAL:HG21	1.92	0.51
1:A:14:THR:OG1	2:B:24:ASP:OD2	2.19	0.51
2:B:153:GLU:OE1	2:B:207:ARG:NH1	2.40	0.51
1:C:57:GLU:OE1	1:C:134:SER:OG	2.28	0.51
3:E:119:LEU:HD23	3:E:119:LEU:C	2.37	0.51
2:D:76:GLY:N	4:G:31:ASN:O	2.43	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:VAL:HG21	1:A:165:MET:HE3	1.93	0.50
2:D:264:ILE:CD1	3:E:271:VAL:HG21	2.41	0.50
1:A:110:MET:HE1	2:B:106:VAL:HG23	1.93	0.50
3:E:225:VAL:HG13	3:E:226:TYR:N	2.28	0.48
2:B:196:ARG:CZ	5:O:1:NAG:H81	2.44	0.48
2:B:121:ASP:OD2	2:B:123:THR:OG1	2.31	0.48
2:D:235:LEU:HD11	3:E:301:LEU:HD23	1.96	0.48
2:B:179:GLU:N	2:B:179:GLU:OE1	2.46	0.48
1:A:110:MET:SD	1:A:111:THR:HA	2.54	0.47
1:C:106:VAL:CG2	7:C:501:NAG:H82	2.44	0.47
2:D:251:VAL:HG12	2:D:255:ILE:HD12	1.96	0.47
2:B:272:LEU:HD13	2:B:279:LYS:HE3	1.97	0.47
3:E:94:LYS:O	3:E:95:LEU:HD23	2.14	0.47
1:A:182:GLU:OE1	1:A:182:GLU:N	2.45	0.47
2:D:73:ALA:O	4:G:102:GLY:N	2.46	0.47
2:D:130:ILE:HG22	2:D:132:THR:HG23	1.97	0.47
2:D:75:SER:N	4:G:100:ARG:O	2.48	0.47
2:B:178:VAL:HA	2:B:181:ILE:HD12	1.97	0.47
1:A:290:VAL:HG11	1:A:406:PHE:CE2	2.49	0.47
1:C:290:VAL:HG11	1:C:406:PHE:CE2	2.50	0.46
1:A:65:ARG:HG2	1:A:128:THR:HG23	1.96	0.46
1:A:116:LEU:HD11	1:A:118:ARG:HE	1.80	0.46
2:D:74:TYR:HA	4:G:101:GLY:HA2	1.97	0.46
3:E:102:ILE:HB	3:E:105:ALA:HB2	1.98	0.46
1:A:102:GLY:O	3:E:133:ARG:NH2	2.49	0.46
3:E:197:PHE:CD1	3:E:215:LEU:HD13	2.50	0.46
9:A:504:PX6:O2	9:A:504:PX6:H5	2.14	0.46
3:E:177:GLU:O	3:E:180:HIS:NE2	2.49	0.46
3:E:312:ASP:OD1	3:E:313:SER:N	2.49	0.46
1:C:20:LEU:HD21	1:C:75:TYR:HB3	1.98	0.45
2:B:258:VAL:HG22	2:B:292:VAL:HG23	1.97	0.45
2:D:420:ALA:O	2:D:421:ALA:HB3	2.17	0.44
2:B:190:GLU:OE2	2:B:192:ARG:NE	2.46	0.44
2:D:178:VAL:HA	2:D:181:ILE:HD12	1.98	0.44
2:D:249:ALA:HB1	3:E:256:VAL:HG21	1.99	0.44
1:C:110:MET:HE1	2:D:106:VAL:CG2	2.47	0.44
2:D:71:ARG:HA	4:G:103:SER:HB3	2.00	0.44
1:C:194:GLN:OE1	1:C:194:GLN:N	2.51	0.44
3:E:163:TYR:HB2	3:E:168:ILE:HD12	1.99	0.44
2:B:166:PHE:HB3	2:B:208:LEU:HD11	1.99	0.44
3:E:256:VAL:O	3:E:260:ILE:HG12	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:290:VAL:HG11	1:C:406:PHE:CZ	2.53	0.44
2:D:253:LEU:CD2	3:E:260:ILE:HG21	2.43	0.44
1:C:290:VAL:HG21	1:C:406:PHE:CE1	2.53	0.43
2:D:305:ILE:HD12	2:D:420:ALA:HB3	2.00	0.43
1:A:11:GLU:N	1:A:11:GLU:OE1	2.52	0.43
2:B:9:MET:HE2	1:C:32:PHE:H	1.83	0.43
1:A:205:THR:O	3:E:121:ARG:NH1	2.51	0.43
3:E:20:LEU:O	3:E:24:ILE:HD12	2.19	0.43
3:E:32:ARG:NH1	3:E:34:GLY:O	2.52	0.43
4:F:36:TRP:HE1	4:F:79:VAL:HG12	1.83	0.43
1:A:283:ALA:HB2	1:A:414:TYR:CZ	2.54	0.42
4:G:36:TRP:HE1	4:G:79:VAL:HG12	1.83	0.42
1:A:55:GLU:OE1	1:A:55:GLU:N	2.52	0.42
1:A:230:ILE:HB	1:A:231:PRO:HD3	2.01	0.42
4:G:98:LEU:O	4:G:99:HIS:HB3	2.20	0.42
4:F:98:LEU:O	4:F:99:HIS:HB3	2.20	0.42
2:D:15:THR:HG21	2:D:77:ILE:CD1	2.50	0.42
2:D:73:ALA:HB3	4:G:102:GLY:O	2.20	0.42
2:D:146:ASP:OD2	2:D:148:GLN:NE2	2.48	0.42
1:A:49:GLY:HA3	1:A:59:THR:HG23	2.02	0.42
2:D:74:TYR:HA	4:G:100:ARG:O	2.20	0.42
2:B:302:VAL:HG13	2:B:420:ALA:HB1	2.02	0.41
1:C:120:THR:OG1	1:C:122:ASN:OD1	2.38	0.41
1:C:259:THR:HG21	2:D:255:ILE:HG21	2.02	0.41
2:D:280:ALA:HB2	2:D:446:TYR:CE1	2.55	0.41
4:G:40:ALA:HB2	4:G:92:ALA:HB2	2.02	0.41
1:A:147:ASP:OD1	1:A:147:ASP:N	2.52	0.41
3:E:192:ILE:HD13	3:E:219:LEU:HD23	2.03	0.41
1:A:307:TYR:CE2	8:A:503:D10:H13	2.56	0.41
2:D:251:VAL:CG1	2:D:255:ILE:HD12	2.51	0.41
2:B:143:TYR:CE2	2:B:272:LEU:HD11	2.56	0.41
1:A:170:THR:HG23	1:A:171:LYS:N	2.36	0.41
1:C:106:VAL:HG23	7:C:501:NAG:H82	2.02	0.41
2:D:183:LEU:HD13	2:D:186:PHE:HB2	2.03	0.41
4:F:40:ALA:HB2	4:F:92:ALA:HB2	2.02	0.41
1:A:73:LEU:HD12	1:A:125:ILE:HD11	2.03	0.41
2:D:120:PRO:O	4:G:33:HIS:CE1	2.69	0.40
2:D:13:LYS:O	2:D:17:ASP:CG	2.65	0.40
2:D:110:THR:O	3:E:110:PHE:HZ	2.05	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	333/364 (92%)	332 (100%)	1 (0%)	0	100	100
1	C	333/364 (92%)	329 (99%)	4 (1%)	0	100	100
2	B	328/451 (73%)	321 (98%)	7 (2%)	0	100	100
2	D	327/451 (72%)	318 (97%)	9 (3%)	0	100	100
3	E	326/430 (76%)	316 (97%)	10 (3%)	0	100	100
4	F	121/123 (98%)	117 (97%)	4 (3%)	0	100	100
4	G	121/123 (98%)	117 (97%)	4 (3%)	0	100	100
All	All	1889/2306 (82%)	1850 (98%)	39 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	301/328 (92%)	301 (100%)	0	100	100
1	C	301/328 (92%)	301 (100%)	0	100	100
2	B	296/393 (75%)	296 (100%)	0	100	100
2	D	296/393 (75%)	294 (99%)	2 (1%)	76	80
3	E	279/363 (77%)	279 (100%)	0	100	100
4	F	99/99 (100%)	99 (100%)	0	100	100
4	G	99/99 (100%)	99 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1671/2003 (83%)	1669 (100%)	2 (0%)	87 91

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	14	GLU
2	D	21	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	41	ASN
2	B	191	HIS
2	D	41	ASN
2	D	85	ASN
2	D	303	ASN
3	E	222	ASN
3	E	308	HIS
4	F	33	HIS
4	G	33	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

5 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	O	1	5	14,14,15	0.43	0	17,19,21	0.79	0
5	BMA	O	2	5	11,11,12	0.21	0	15,15,17	0.64	0
6	NAG	P	1	2,6	14,14,15	0.49	0	17,19,21	1.09	1 (5%)
6	NAG	P	2	6	14,14,15	0.46	0	17,19,21	0.95	1 (5%)
6	BMA	P	3	6	11,11,12	0.20	0	15,15,17	1.09	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	O	1	5	-	4/6/23/26	0/1/1/1
5	BMA	O	2	5	-	0/2/19/22	0/1/1/1
6	NAG	P	1	2,6	-	2/6/23/26	0/1/1/1
6	NAG	P	2	6	-	4/6/23/26	0/1/1/1
6	BMA	P	3	6	-	1/2/19/22	1/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	P	1	NAG	C1-C2-N2	3.86	117.08	110.49
6	P	3	BMA	C1-O5-C5	3.27	116.62	112.19
6	P	2	NAG	C1-O5-C5	-2.25	109.14	112.19

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	O	1	NAG	C8-C7-N2-C2
5	O	1	NAG	O7-C7-N2-C2
6	P	1	NAG	C8-C7-N2-C2
6	P	1	NAG	O7-C7-N2-C2
6	P	2	NAG	C8-C7-N2-C2
6	P	2	NAG	O7-C7-N2-C2
5	O	1	NAG	O5-C5-C6-O6
6	P	2	NAG	O5-C5-C6-O6
6	P	3	BMA	O5-C5-C6-O6
5	O	1	NAG	C4-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
6	P	2	NAG	C4-C5-C6-O6

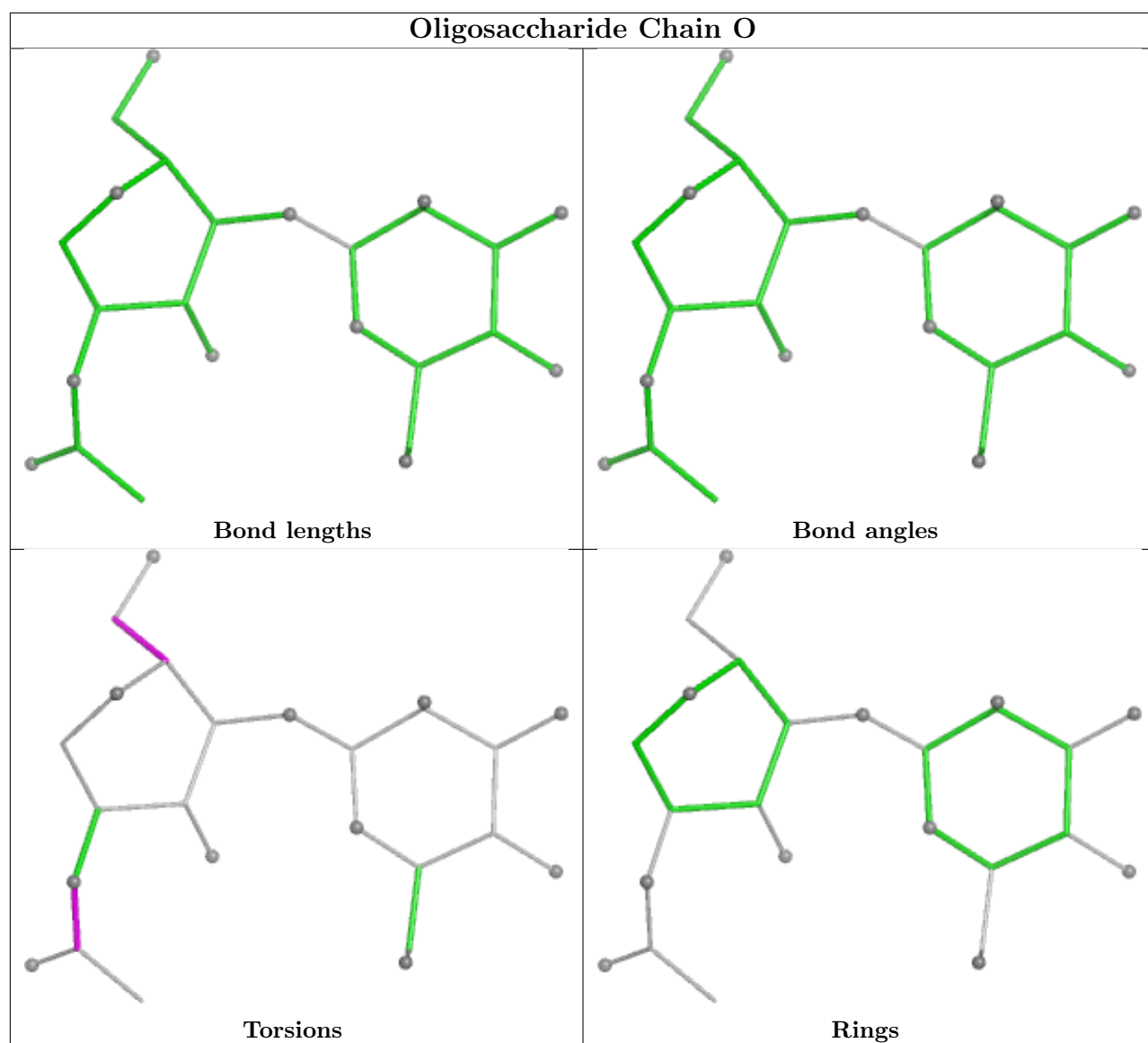
All (1) ring outliers are listed below:

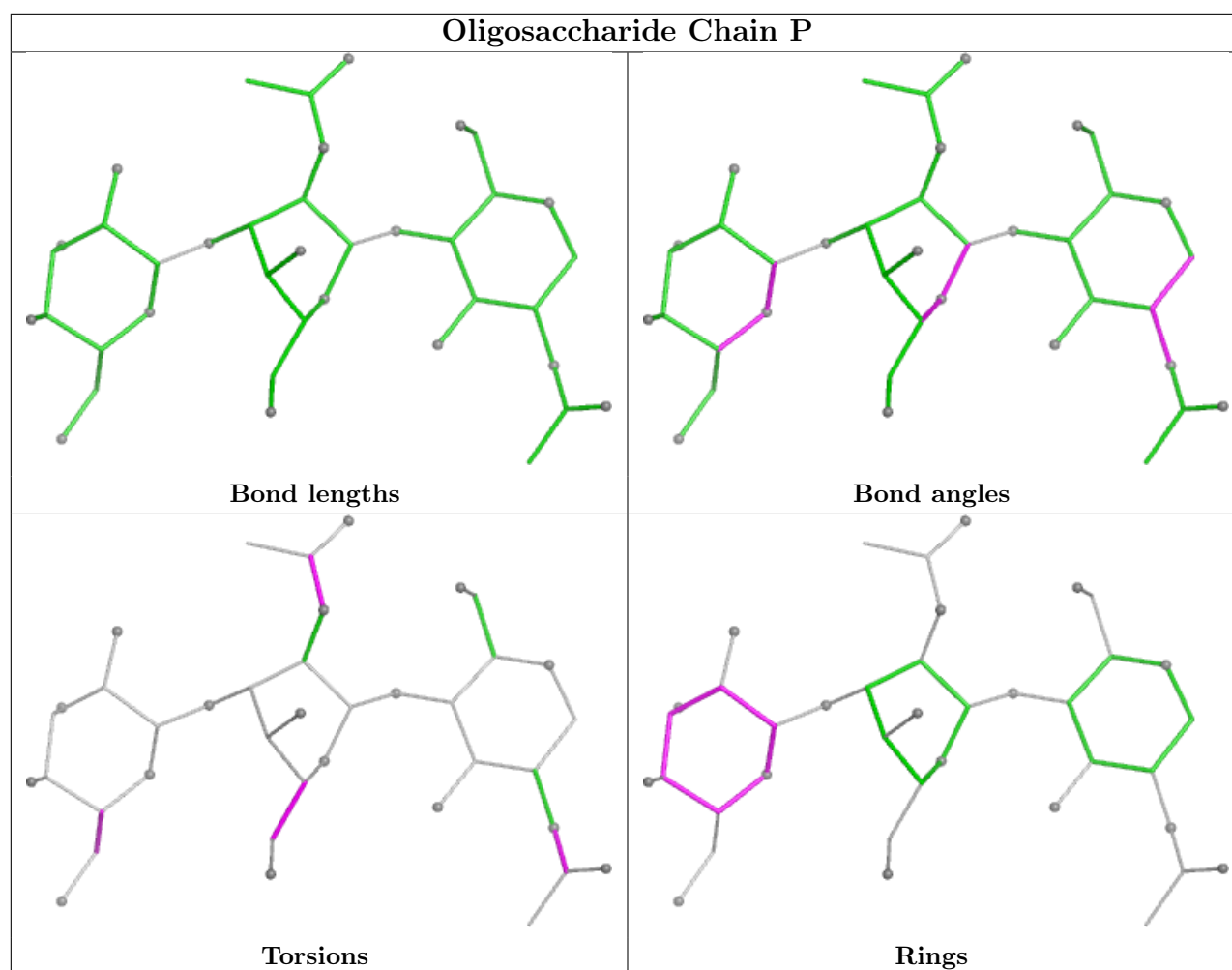
Mol	Chain	Res	Type	Atoms
6	P	3	BMA	C1-C2-C3-C4-C5-O5

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	O	1	NAG	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

15 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
9	PX6	A	504	-	43,43,43	0.28	0	47,48,48	0.33	0
8	D10	B	505	-	9,9,9	0.16	0	8,8,8	0.12	0
7	NAG	B	504	2	14,14,15	0.22	0	17,19,21	0.45	0
7	NAG	B	503	2	14,14,15	0.23	0	17,19,21	0.49	0
7	NAG	A	502	1	14,14,15	0.22	0	17,19,21	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	D10	A	503	-	9,9,9	0.16	0	8,8,8	0.11	0
10	ABU	D	501	-	6,6,6	0.77	0	6,6,6	0.10	0
7	NAG	A	501	1	14,14,15	0.23	0	17,19,21	0.59	0
8	D10	D	503	-	9,9,9	0.17	0	8,8,8	0.17	0
7	NAG	E	501	3	14,14,15	0.23	0	17,19,21	0.47	0
7	NAG	D	502	2	14,14,15	0.22	0	17,19,21	0.45	0
7	NAG	B	502	2	14,14,15	0.38	0	17,19,21	1.02	2 (11%)
8	D10	A	505	-	9,9,9	0.17	0	8,8,8	0.10	0
10	ABU	B	501	-	6,6,6	0.77	0	6,6,6	0.12	0
7	NAG	C	501	1	14,14,15	0.23	0	17,19,21	0.89	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	PX6	A	504	-	-	6/45/45/45	-
8	D10	B	505	-	-	0/7/7/7	-
7	NAG	B	504	2	-	0/6/23/26	0/1/1/1
7	NAG	B	503	2	-	0/6/23/26	0/1/1/1
7	NAG	A	502	1	-	0/6/23/26	0/1/1/1
8	D10	A	503	-	-	0/7/7/7	-
10	ABU	D	501	-	-	2/4/4/4	-
7	NAG	A	501	1	-	0/6/23/26	0/1/1/1
8	D10	D	503	-	-	0/7/7/7	-
7	NAG	E	501	3	-	0/6/23/26	0/1/1/1
7	NAG	D	502	2	-	0/6/23/26	0/1/1/1
7	NAG	B	502	2	-	4/6/23/26	0/1/1/1
8	D10	A	505	-	-	0/7/7/7	-
10	ABU	B	501	-	-	2/4/4/4	-
7	NAG	C	501	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	501	NAG	C1-O5-C5	3.32	116.69	112.19
7	B	502	NAG	C1-O5-C5	2.49	115.57	112.19
7	B	502	NAG	C2-N2-C7	2.24	126.09	122.90

There are no chirality outliers.

All (16) torsion outliers are listed below:

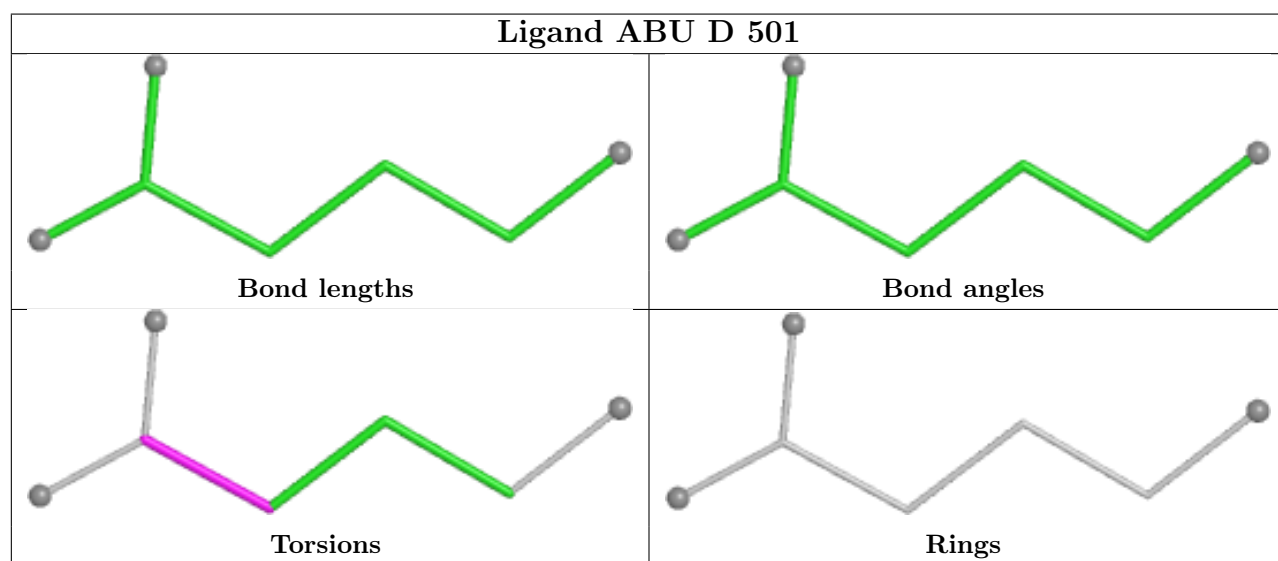
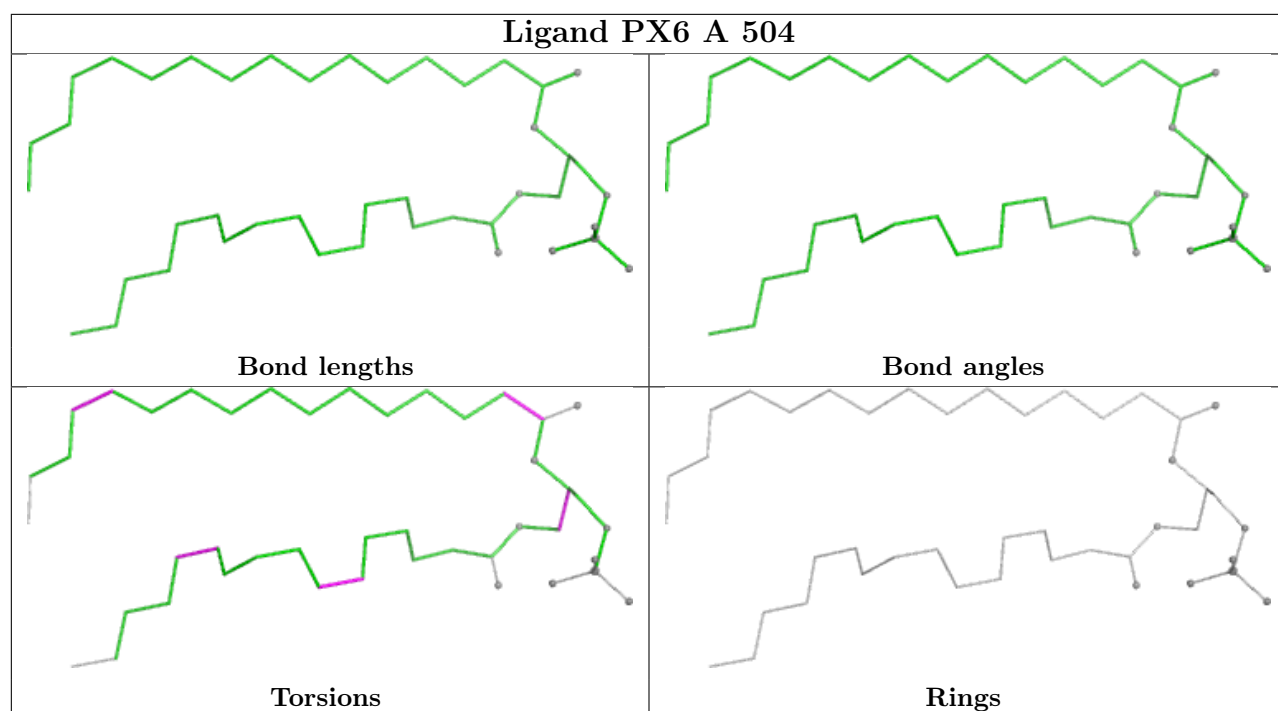
Mol	Chain	Res	Type	Atoms
7	B	502	NAG	C3-C2-N2-C7
7	B	502	NAG	C8-C7-N2-C2
7	B	502	NAG	O7-C7-N2-C2
9	A	504	PX6	O7-C2-C3-O5
9	A	504	PX6	C11-C10-C9-C8
7	C	501	NAG	O5-C5-C6-O6
7	B	502	NAG	O5-C5-C6-O6
9	A	504	PX6	C1-C2-C3-O5
7	C	501	NAG	C3-C2-N2-C7
9	A	504	PX6	O7-C20-C21-C22
9	A	504	PX6	C13-C14-C15-C16
10	B	501	ABU	OXT-C-CG-CB
10	D	501	ABU	OXT-C-CG-CB
10	D	501	ABU	O-C-CG-CB
10	B	501	ABU	O-C-CG-CB
9	A	504	PX6	C30-C31-C32-C33

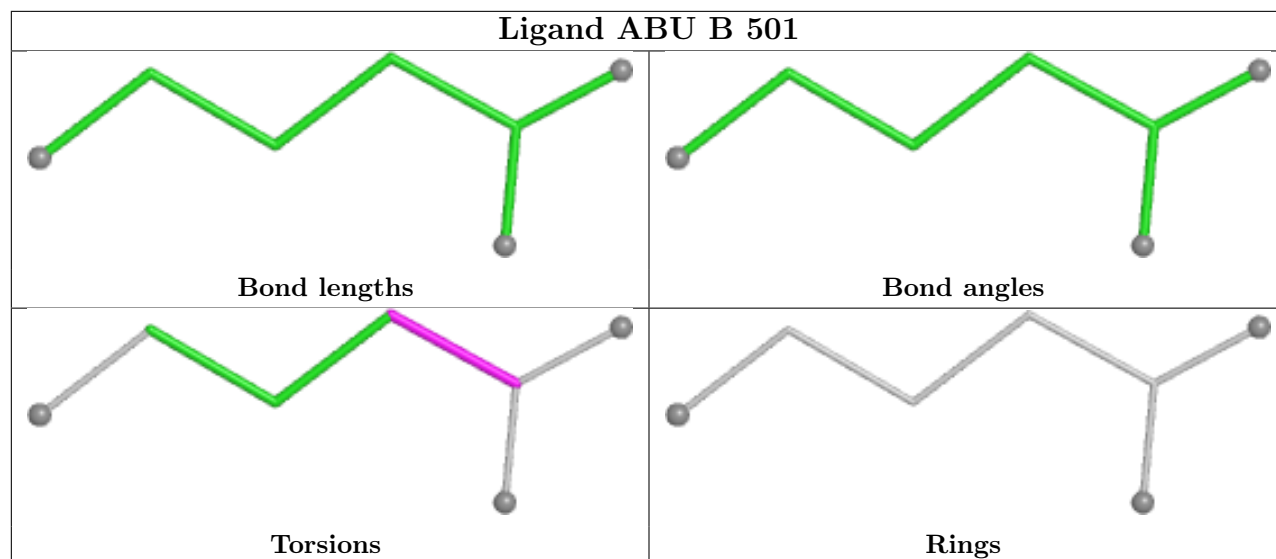
There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	504	PX6	3	0
8	A	503	D10	1	0
7	B	502	NAG	4	0
7	C	501	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.

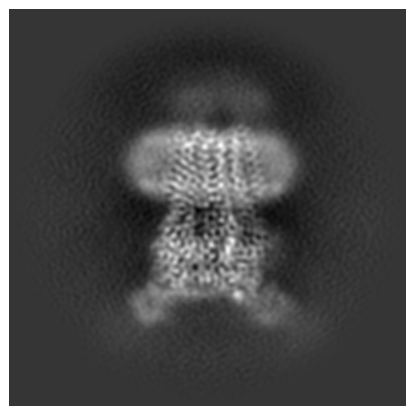
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55007. These allow visual inspection of the internal detail of the map and identification of artifacts.

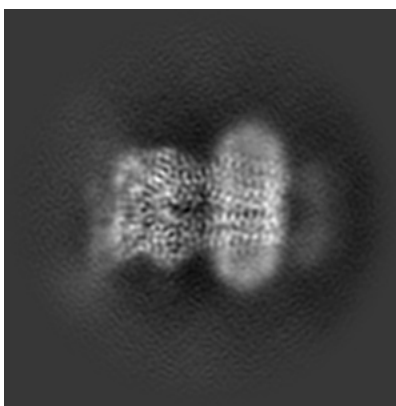
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

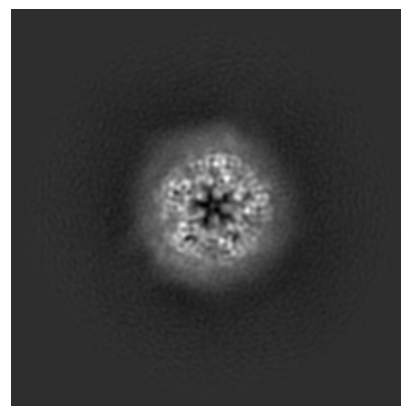
6.1.1 Primary map



X

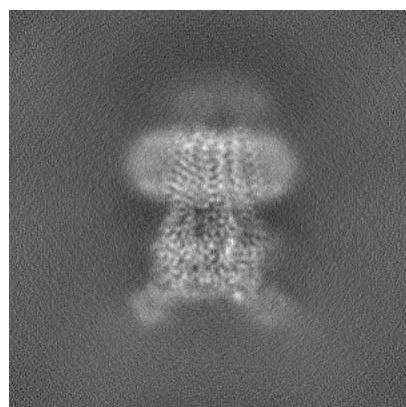


Y

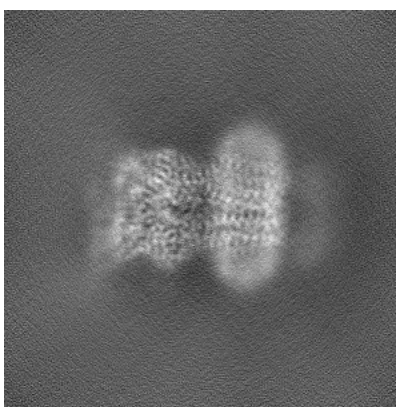


Z

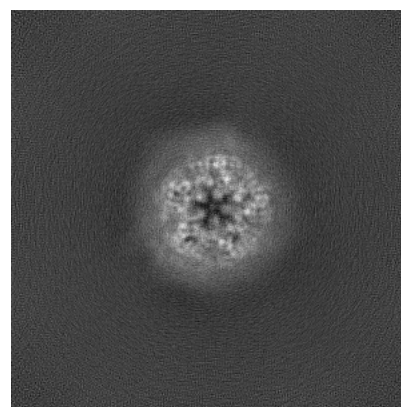
6.1.2 Raw map



X



Y

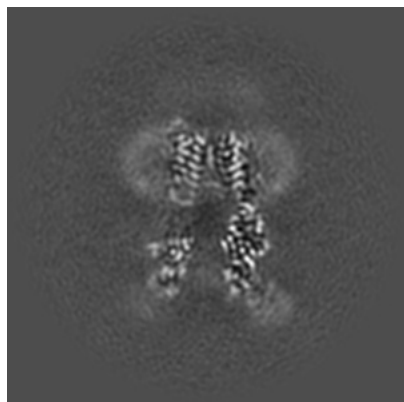


Z

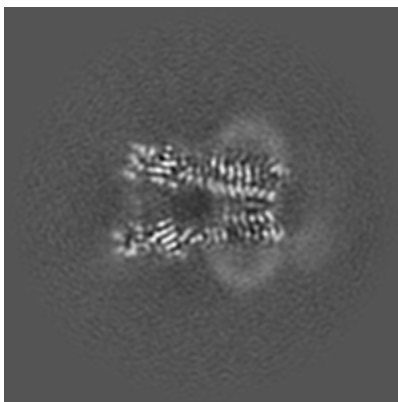
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

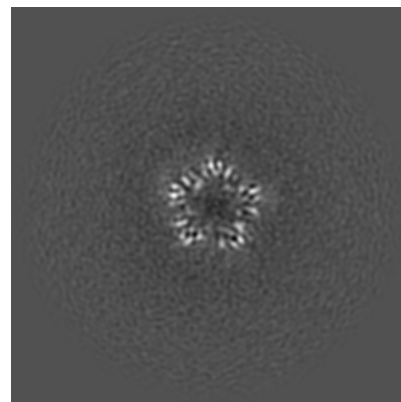
6.2.1 Primary map



X Index: 192

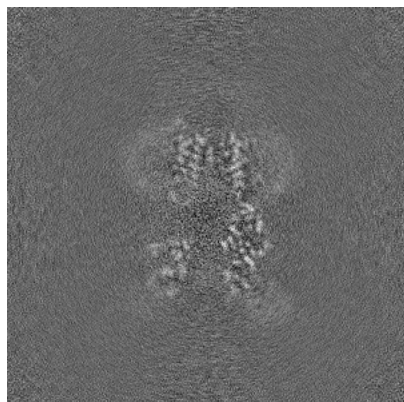


Y Index: 192

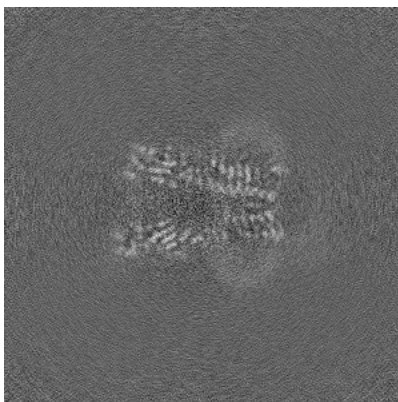


Z Index: 192

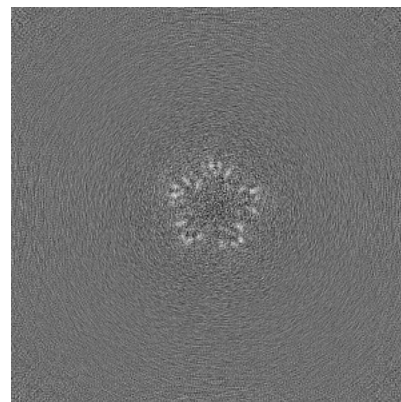
6.2.2 Raw map



X Index: 192



Y Index: 192

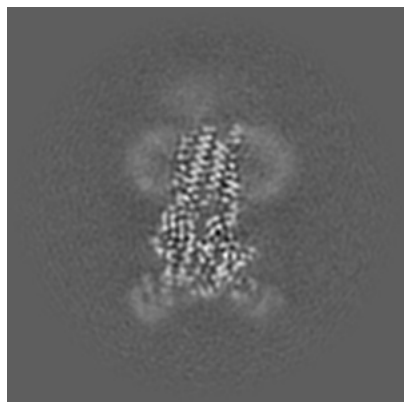


Z Index: 192

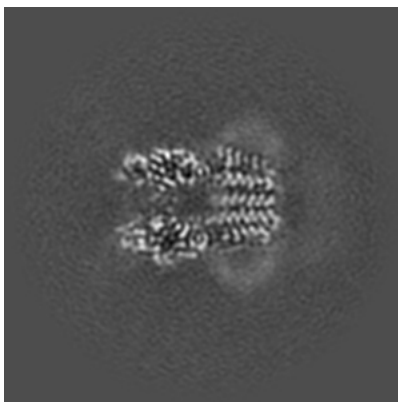
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

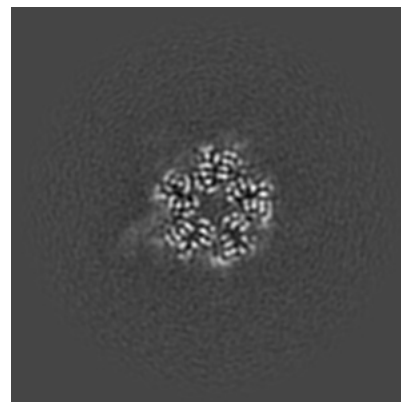
6.3.1 Primary map



X Index: 167

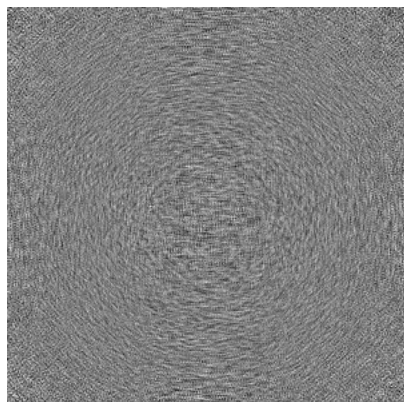


Y Index: 204

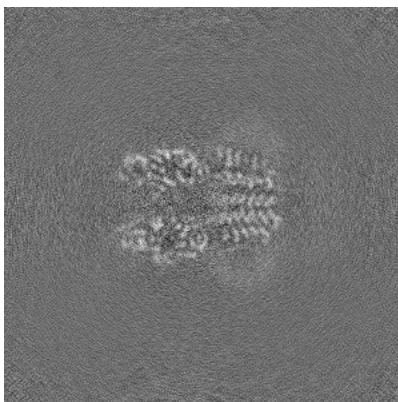


Z Index: 156

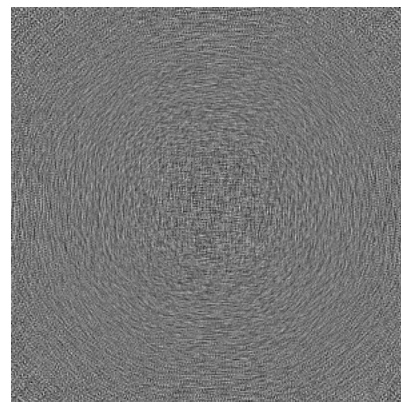
6.3.2 Raw map



X Index: 1



Y Index: 204

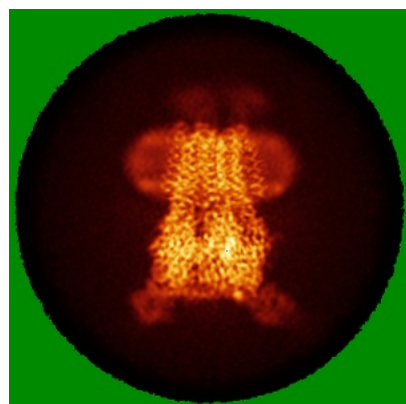


Z Index: 0

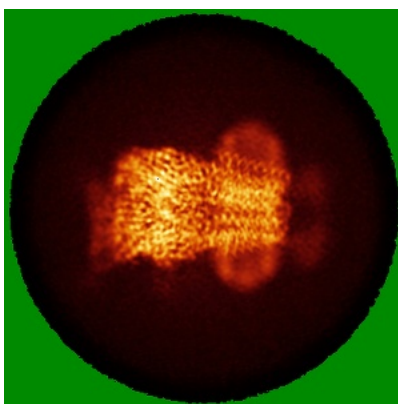
The images above show the largest variance slices of the map in three orthogonal directions.

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

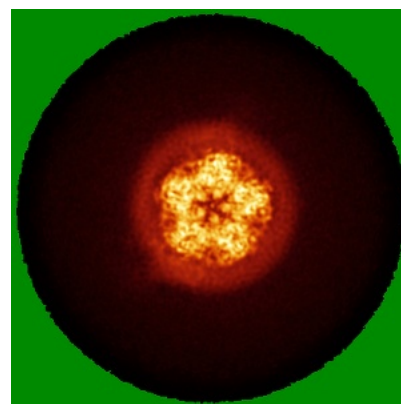
6.4.1 Primary map



X

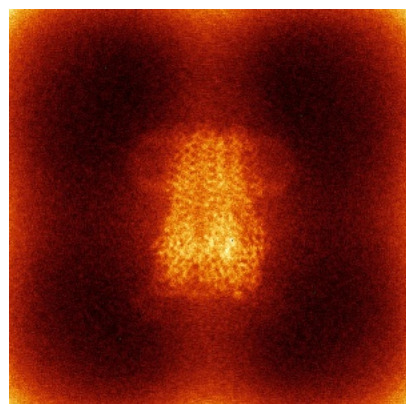


Y

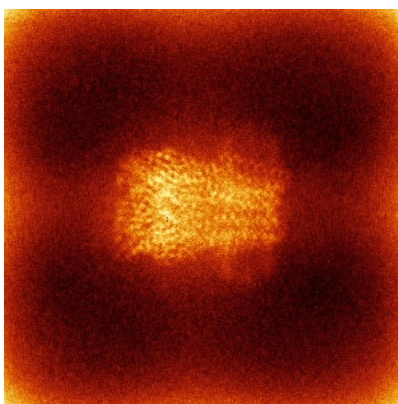


Z

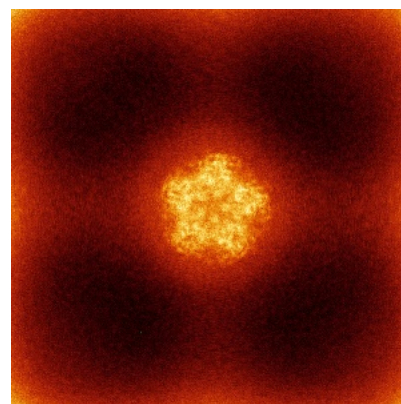
6.4.2 Raw map



X



Y

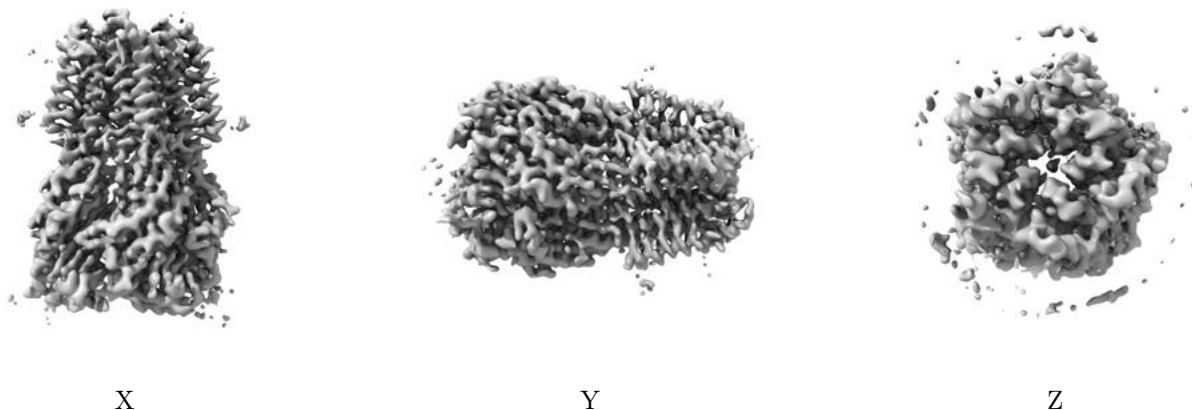


Z

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.

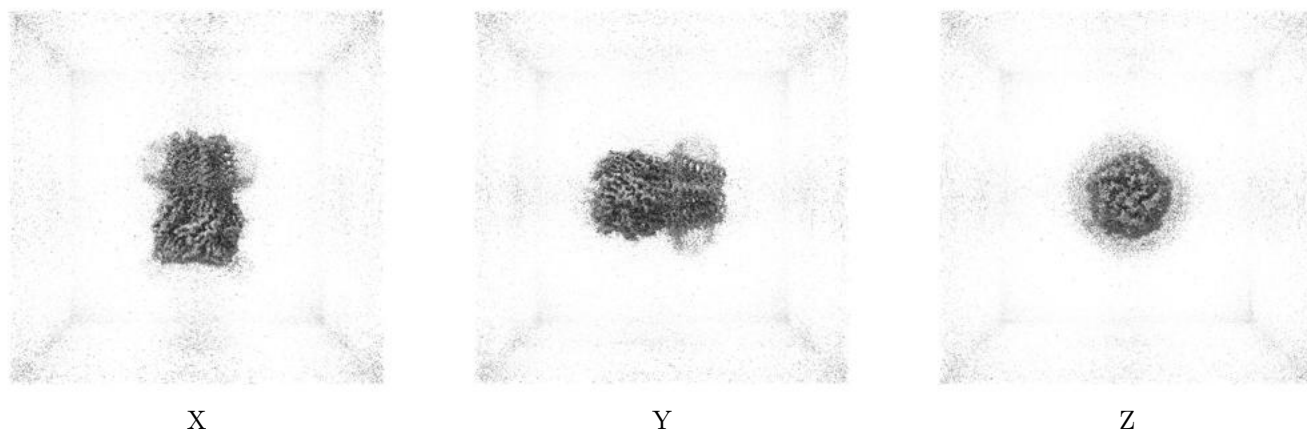
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0559. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

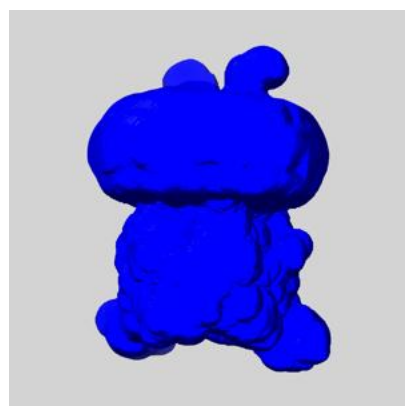
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

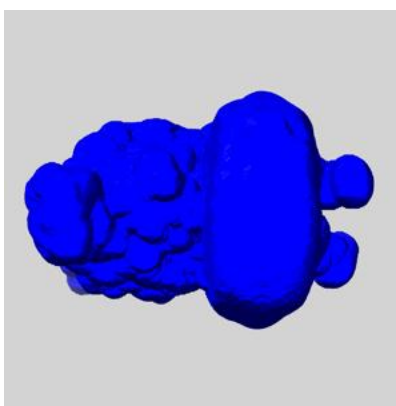
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

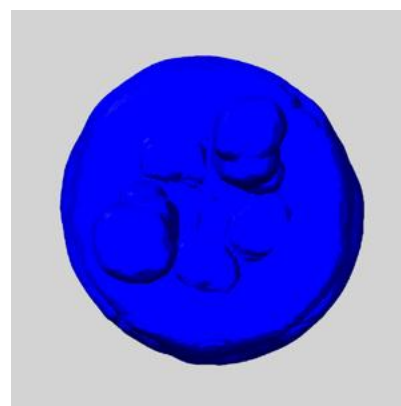
6.6.1 emd_55007_msk_1.map [i](#)



X



Y

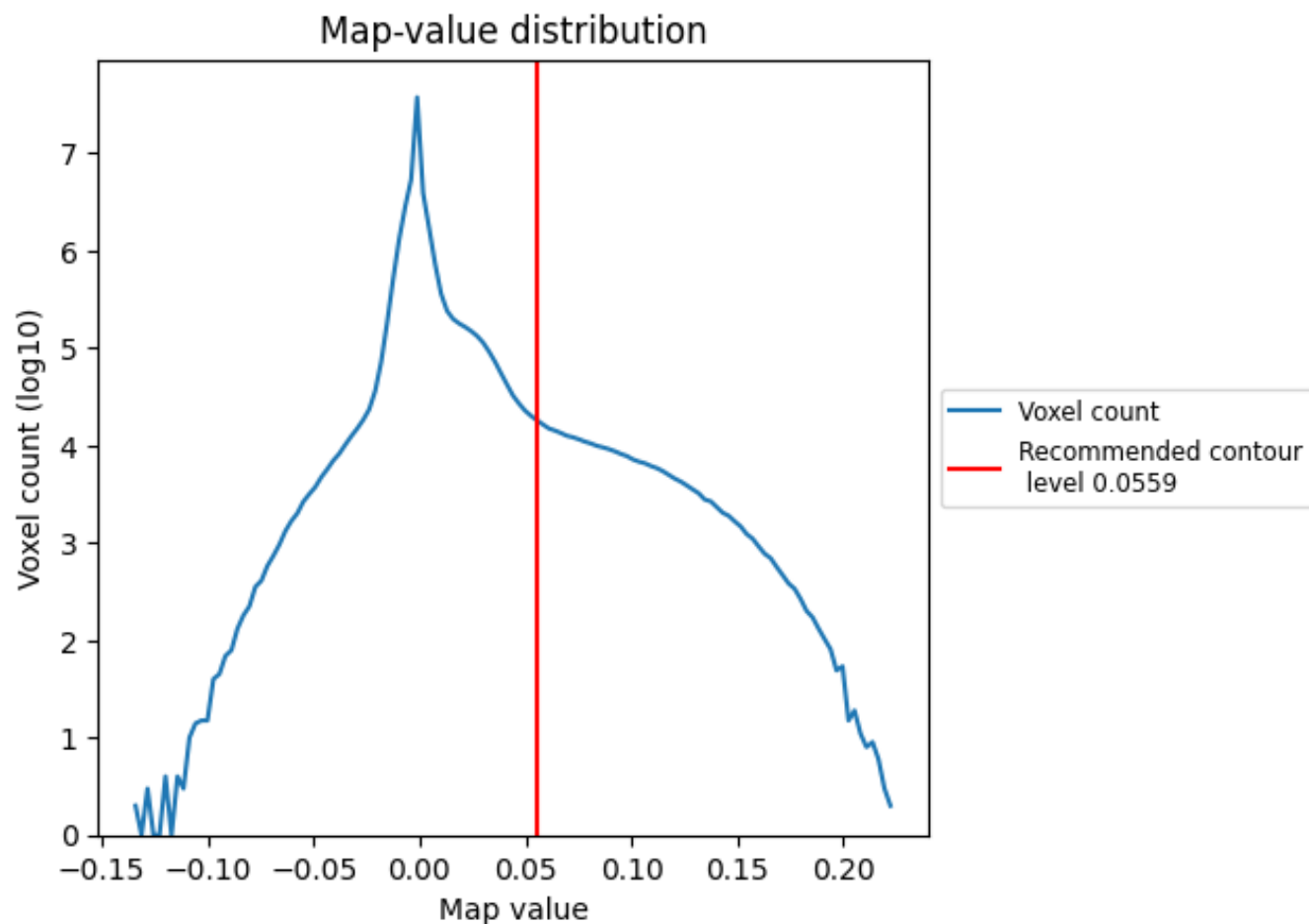


Z

7 Map analysis [i](#)

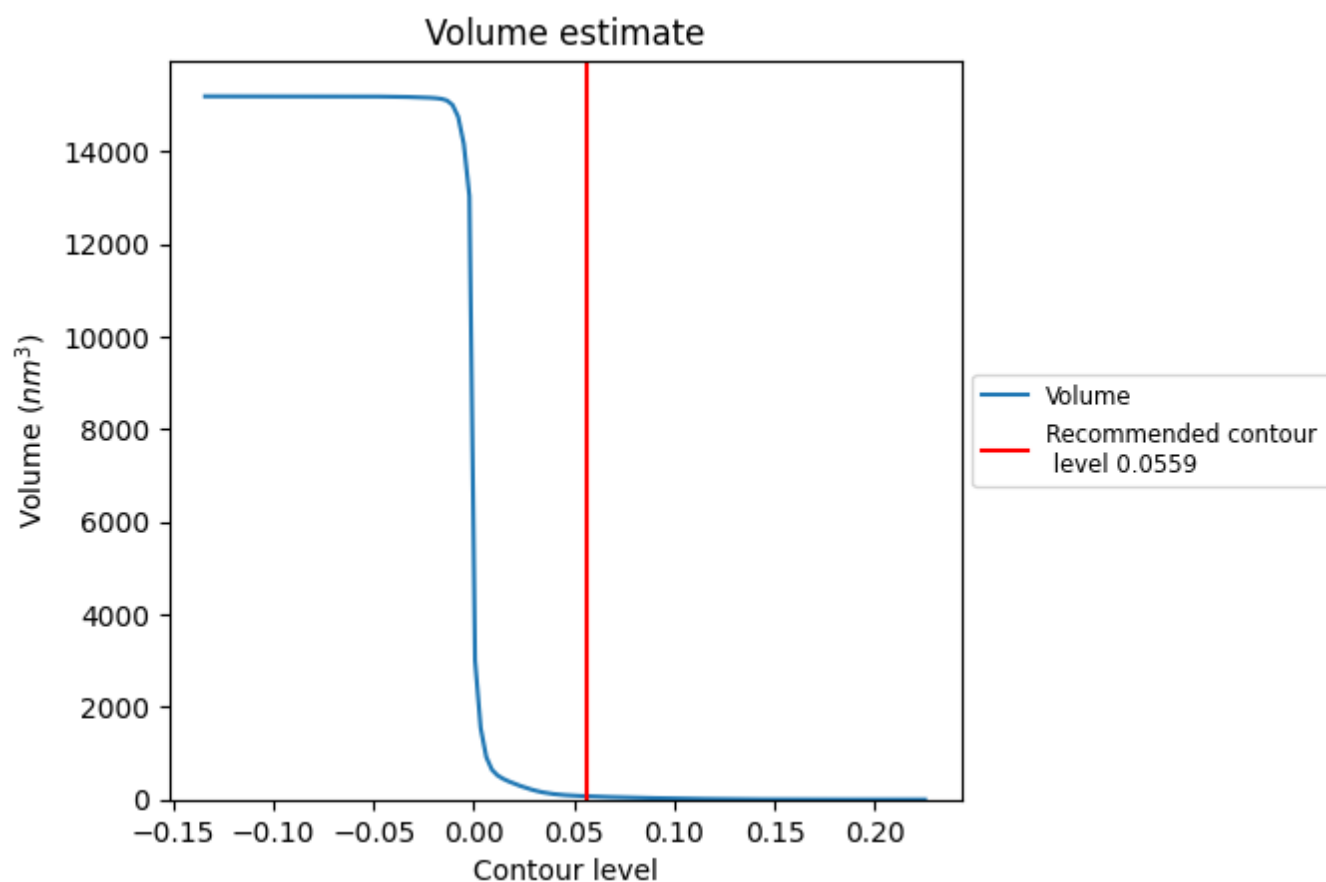
This section contains the results of statistical analysis of the map.

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.

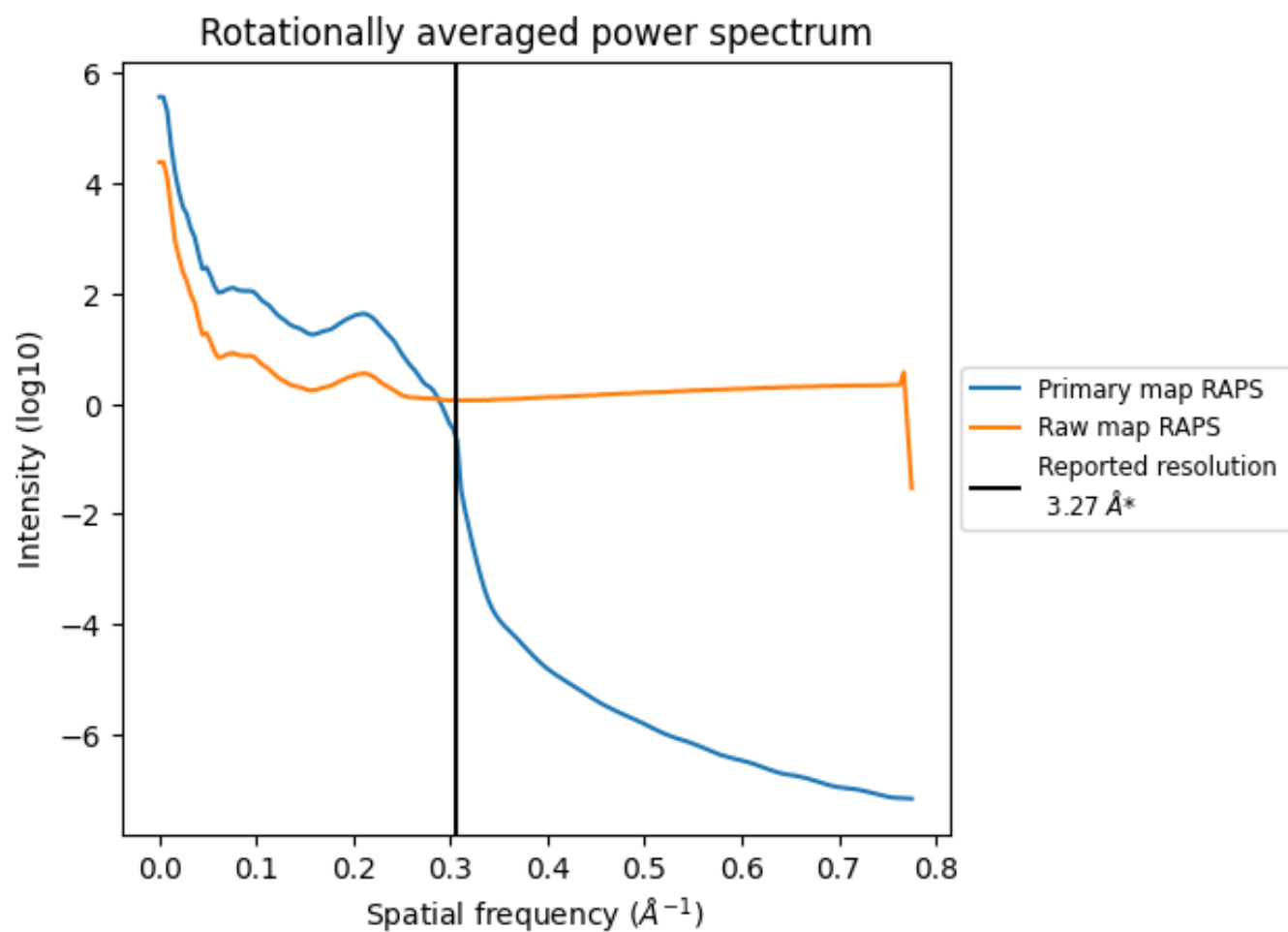
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 73 nm³; this corresponds to an approximate mass of 66 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 ⓘ

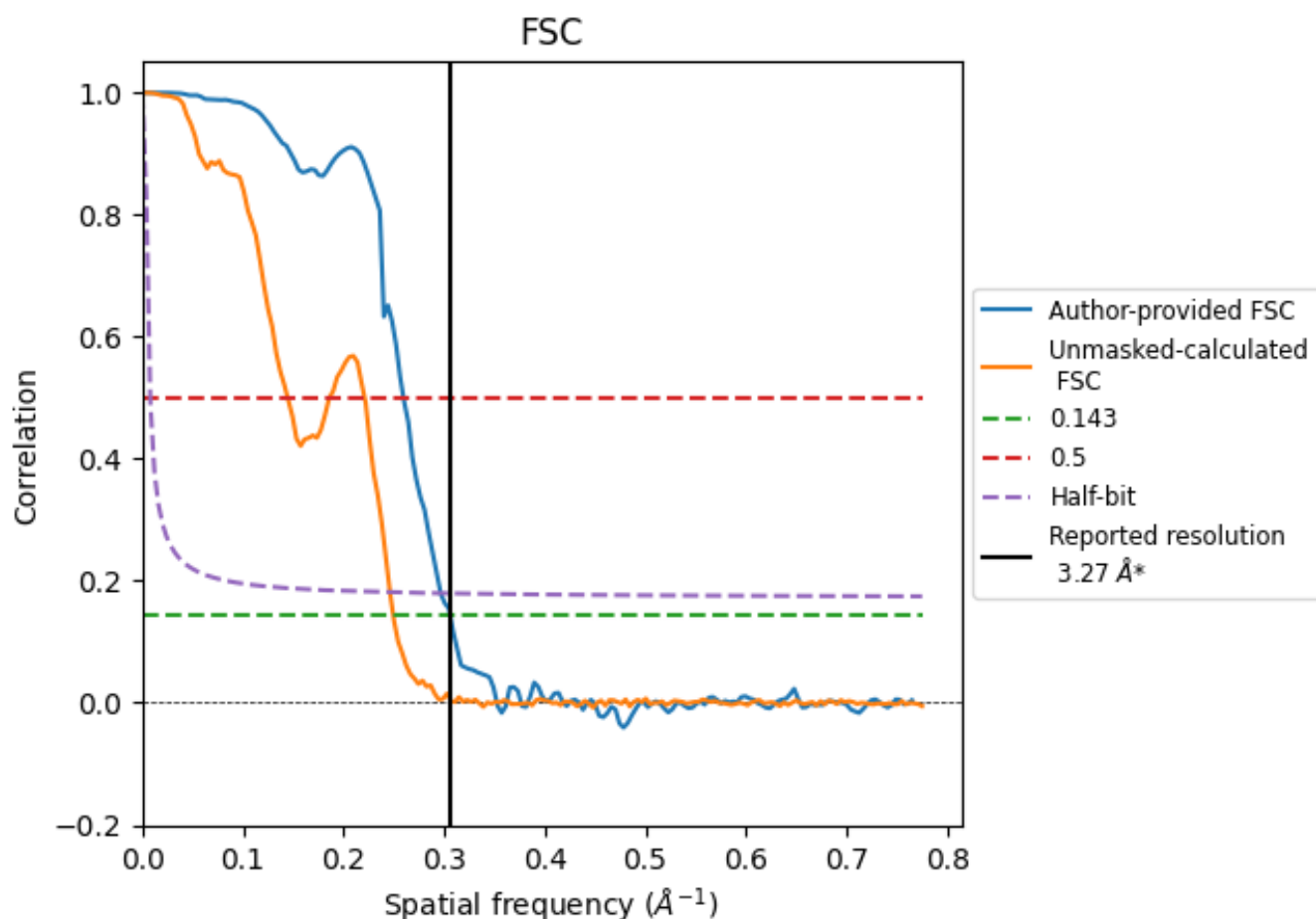


*Reported resolution corresponds to spatial frequency of 0.306 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.306 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

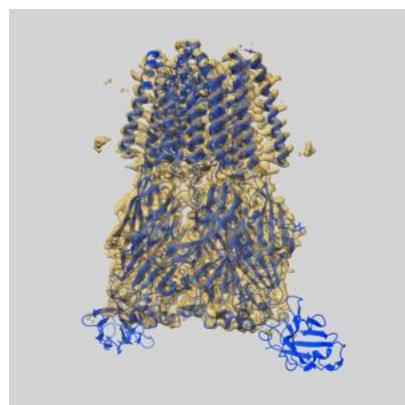
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.27	-	-
Author-provided FSC curve	3.27	3.85	3.37
Unmasked-calculated*	4.01	6.92	4.06

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.01 differs from the reported value 3.27 by more than 10 %

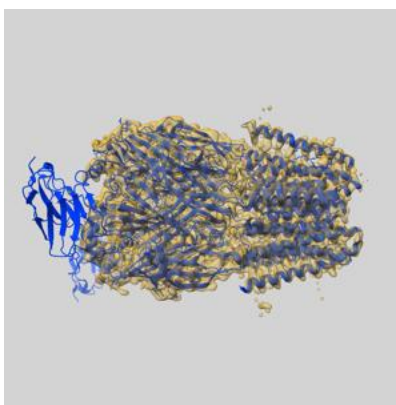
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-55007 and PDB model 9SLM. Per-residue inclusion information can be found in section 3 on page 20.

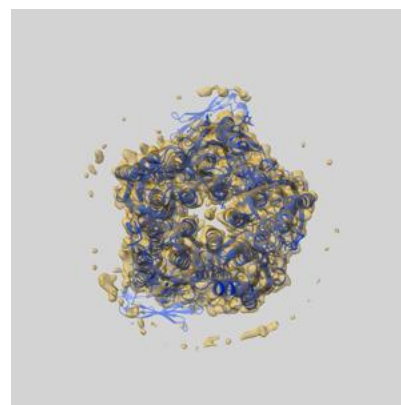
9.1 Map-model overlay [i](#)



X



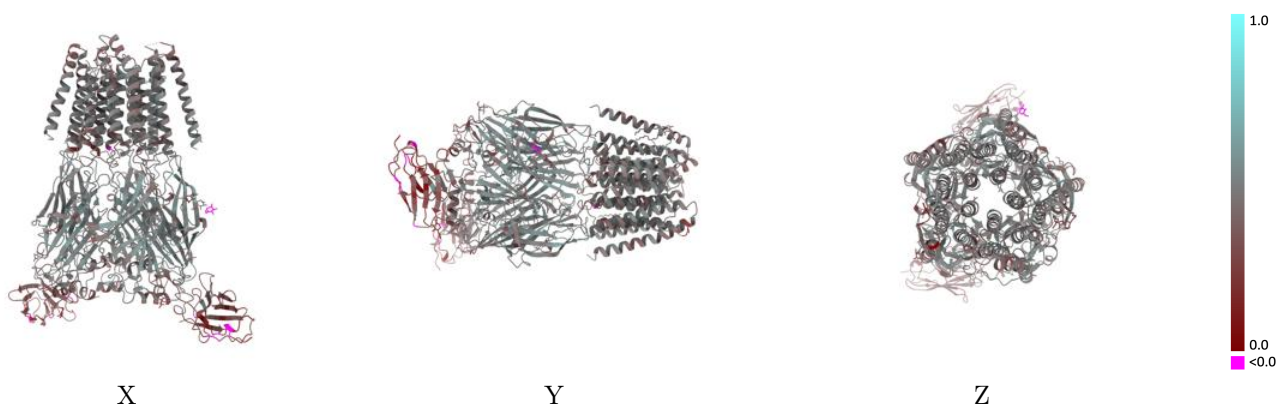
Y



Z

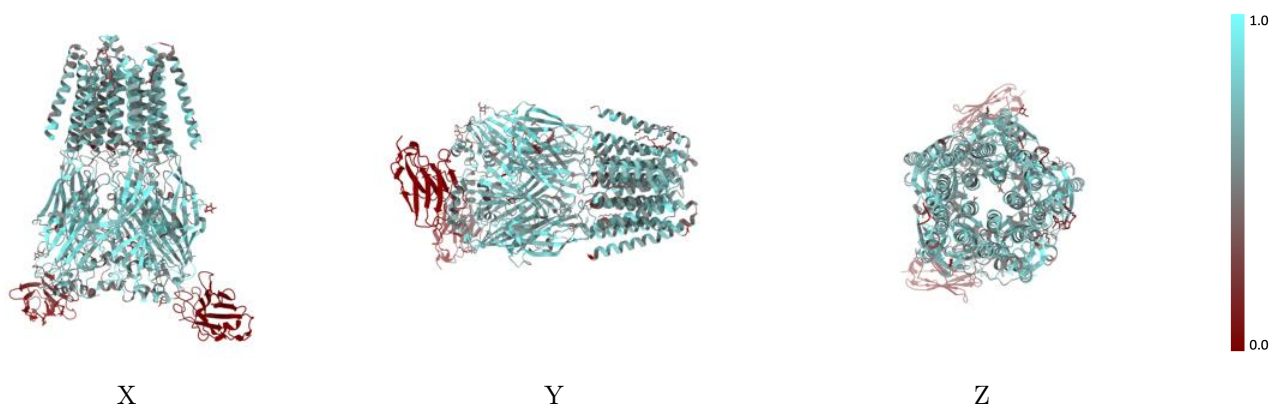
The images above show the 3D surface view of the map at the recommended contour level 0.0559 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



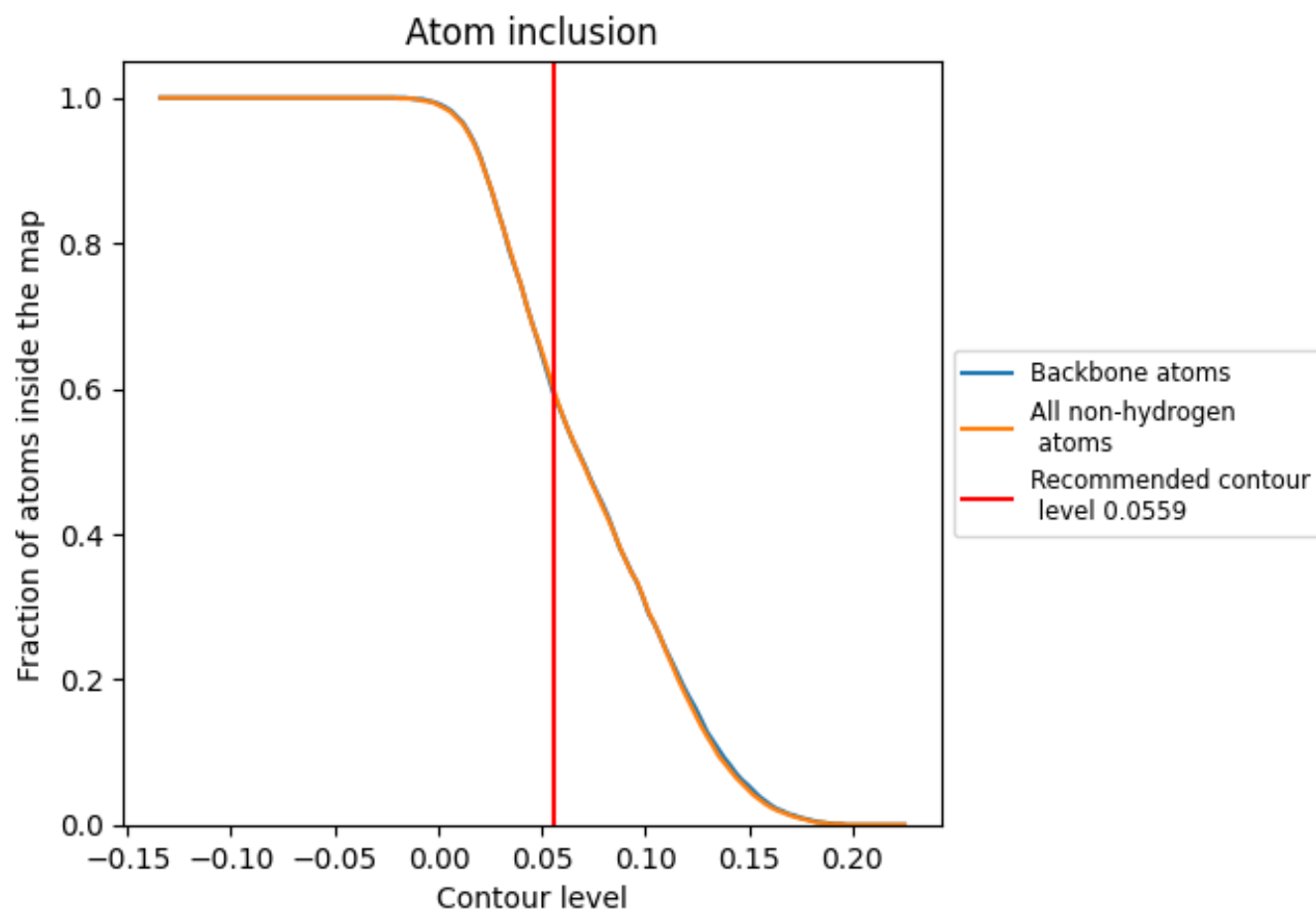
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0559).

9.4 Atom inclusion [i](#)



At the recommended contour level, 59% of all backbone atoms, 60% 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 (0.0559) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5970	0.4350
A	0.6870	0.4640
B	0.6890	0.4720
C	0.6720	0.4590
D	0.6750	0.4520
E	0.6810	0.4570
F	0.0210	0.2610
G	0.0390	0.2500
O	0.4000	0.1420
P	0.5900	0.3550

1.0

0.0

<0.0