



## Full wwPDB EM Validation Report ⓘ

Nov 3, 2024 – 05:26 AM EST

PDB ID : 7KEW  
EMDB ID : EMD-22841  
Title : Bundibugyo virus GP (mucin deleted) bound to antibody Fab BDBV-43  
Authors : Murin, C.D.; Ward, A.B.  
Deposited on : 2020-10-13  
Resolution : 4.16 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.dev113  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

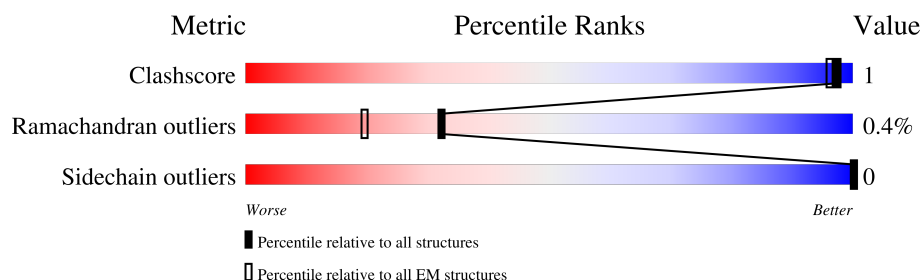
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*









The reported resolution of this entry is 4.16 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	343	 62% 34%
1	B	343	 5% 62% 34%
1	C	343	 5% 62% 34%
2	G	246	 7% 48% 50%
2	H	246	 7% 48% 50%
2	I	246	 7% 48% 50%
3	J	232	 16% 43% 54%
3	K	232	 16% 44% 54%

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Mol	Chain	Length	Quality of chain
3	L	232	
4	D	177	
4	E	177	
4	F	177	
5	M	2	
5	N	2	
5	Q	2	
6	O	5	
6	P	5	
6	R	5	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 13173 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 Spike glycoprotein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	227	Total	C	N	O	S	0	0
			1785	1139	304	336	6		
1	B	227	Total	C	N	O	S	0	0
			1785	1139	304	336	6		
1	C	227	Total	C	N	O	S	0	0
			1785	1139	304	336	6		

There are 93 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	313	ILE	-	expression tag	UNP A0A510C2V9
A	314	ASP	-	expression tag	UNP A0A510C2V9
A	315	ILE	-	expression tag	UNP A0A510C2V9
A	316	SER	-	expression tag	UNP A0A510C2V9
A	317	GLU	-	expression tag	UNP A0A510C2V9
A	318	SER	-	expression tag	UNP A0A510C2V9
A	319	THR	-	expression tag	UNP A0A510C2V9
A	320	GLU	-	expression tag	UNP A0A510C2V9
A	321	PRO	-	expression tag	UNP A0A510C2V9
A	322	GLY	-	expression tag	UNP A0A510C2V9
A	323	PRO	-	expression tag	UNP A0A510C2V9
A	324	LEU	-	expression tag	UNP A0A510C2V9
A	325	THR	-	expression tag	UNP A0A510C2V9
A	326	ASN	-	expression tag	UNP A0A510C2V9
A	327	THR	-	expression tag	UNP A0A510C2V9
A	328	THR	-	expression tag	UNP A0A510C2V9
A	329	ARG	-	expression tag	UNP A0A510C2V9
A	330	GLY	-	expression tag	UNP A0A510C2V9
A	331	ALA	-	expression tag	UNP A0A510C2V9
A	332	ALA	-	expression tag	UNP A0A510C2V9
A	333	ASN	-	expression tag	UNP A0A510C2V9
A	334	LEU	-	expression tag	UNP A0A510C2V9
A	335	LEU	-	expression tag	UNP A0A510C2V9
A	336	THR	-	expression tag	UNP A0A510C2V9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	337	GLY	-	expression tag	UNP A0A510C2V9
A	338	SER	-	expression tag	UNP A0A510C2V9
A	339	ARG	-	expression tag	UNP A0A510C2V9
A	340	ARG	-	expression tag	UNP A0A510C2V9
A	341	THR	-	expression tag	UNP A0A510C2V9
A	342	ARG	-	expression tag	UNP A0A510C2V9
A	343	ARG	-	expression tag	UNP A0A510C2V9
B	313	ILE	-	expression tag	UNP A0A510C2V9
B	314	ASP	-	expression tag	UNP A0A510C2V9
B	315	ILE	-	expression tag	UNP A0A510C2V9
B	316	SER	-	expression tag	UNP A0A510C2V9
B	317	GLU	-	expression tag	UNP A0A510C2V9
B	318	SER	-	expression tag	UNP A0A510C2V9
B	319	THR	-	expression tag	UNP A0A510C2V9
B	320	GLU	-	expression tag	UNP A0A510C2V9
B	321	PRO	-	expression tag	UNP A0A510C2V9
B	322	GLY	-	expression tag	UNP A0A510C2V9
B	323	PRO	-	expression tag	UNP A0A510C2V9
B	324	LEU	-	expression tag	UNP A0A510C2V9
B	325	THR	-	expression tag	UNP A0A510C2V9
B	326	ASN	-	expression tag	UNP A0A510C2V9
B	327	THR	-	expression tag	UNP A0A510C2V9
B	328	THR	-	expression tag	UNP A0A510C2V9
B	329	ARG	-	expression tag	UNP A0A510C2V9
B	330	GLY	-	expression tag	UNP A0A510C2V9
B	331	ALA	-	expression tag	UNP A0A510C2V9
B	332	ALA	-	expression tag	UNP A0A510C2V9
B	333	ASN	-	expression tag	UNP A0A510C2V9
B	334	LEU	-	expression tag	UNP A0A510C2V9
B	335	LEU	-	expression tag	UNP A0A510C2V9
B	336	THR	-	expression tag	UNP A0A510C2V9
B	337	GLY	-	expression tag	UNP A0A510C2V9
B	338	SER	-	expression tag	UNP A0A510C2V9
B	339	ARG	-	expression tag	UNP A0A510C2V9
B	340	ARG	-	expression tag	UNP A0A510C2V9
B	341	THR	-	expression tag	UNP A0A510C2V9
B	342	ARG	-	expression tag	UNP A0A510C2V9
B	343	ARG	-	expression tag	UNP A0A510C2V9
C	313	ILE	-	expression tag	UNP A0A510C2V9
C	314	ASP	-	expression tag	UNP A0A510C2V9
C	315	ILE	-	expression tag	UNP A0A510C2V9
C	316	SER	-	expression tag	UNP A0A510C2V9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	317	GLU	-	expression tag	UNP A0A510C2V9
C	318	SER	-	expression tag	UNP A0A510C2V9
C	319	THR	-	expression tag	UNP A0A510C2V9
C	320	GLU	-	expression tag	UNP A0A510C2V9
C	321	PRO	-	expression tag	UNP A0A510C2V9
C	322	GLY	-	expression tag	UNP A0A510C2V9
C	323	PRO	-	expression tag	UNP A0A510C2V9
C	324	LEU	-	expression tag	UNP A0A510C2V9
C	325	THR	-	expression tag	UNP A0A510C2V9
C	326	ASN	-	expression tag	UNP A0A510C2V9
C	327	THR	-	expression tag	UNP A0A510C2V9
C	328	THR	-	expression tag	UNP A0A510C2V9
C	329	ARG	-	expression tag	UNP A0A510C2V9
C	330	GLY	-	expression tag	UNP A0A510C2V9
C	331	ALA	-	expression tag	UNP A0A510C2V9
C	332	ALA	-	expression tag	UNP A0A510C2V9
C	333	ASN	-	expression tag	UNP A0A510C2V9
C	334	LEU	-	expression tag	UNP A0A510C2V9
C	335	LEU	-	expression tag	UNP A0A510C2V9
C	336	THR	-	expression tag	UNP A0A510C2V9
C	337	GLY	-	expression tag	UNP A0A510C2V9
C	338	SER	-	expression tag	UNP A0A510C2V9
C	339	ARG	-	expression tag	UNP A0A510C2V9
C	340	ARG	-	expression tag	UNP A0A510C2V9
C	341	THR	-	expression tag	UNP A0A510C2V9
C	342	ARG	-	expression tag	UNP A0A510C2V9
C	343	ARG	-	expression tag	UNP A0A510C2V9

- Molecule 2 is a protein called BDBV-43 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	G	122	Total	C	N	O	S	0	0
			931	583	163	180	5		
2	H	122	Total	C	N	O	S	0	0
			931	583	163	180	5		
2	I	122	Total	C	N	O	S	0	0
			931	583	163	180	5		

- Molecule 3 is a protein called BDBV-43 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	106	Total	C	N	O	S	0	0
			815	514	141	156	4		
3	K	106	Total	C	N	O	S	0	0
			815	514	141	156	4		
3	L	106	Total	C	N	O	S	0	0
			815	514	141	156	4		

- Molecule 4 is a protein called Envelope glycoprotein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	95	Total	C	N	O	S	0	0
			757	481	132	138	6		
4	E	95	Total	C	N	O	S	0	0
			757	481	132	138	6		
4	F	95	Total	C	N	O	S	0	0
			757	481	132	138	6		

There are 114 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	641	VAL	-	expression tag	UNP B8XCNO
D	642	GLU	-	expression tag	UNP B8XCNO
D	643	VAL	-	expression tag	UNP B8XCNO
D	644	ASP	-	expression tag	UNP B8XCNO
D	645	ASP	-	expression tag	UNP B8XCNO
D	646	ASP	-	expression tag	UNP B8XCNO
D	647	ASP	-	expression tag	UNP B8XCNO
D	648	LYS	-	expression tag	UNP B8XCNO
D	649	ALA	-	expression tag	UNP B8XCNO
D	650	GLY	-	expression tag	UNP B8XCNO
D	651	TRP	-	expression tag	UNP B8XCNO
D	652	SER	-	expression tag	UNP B8XCNO
D	653	HIS	-	expression tag	UNP B8XCNO
D	654	PRO	-	expression tag	UNP B8XCNO
D	655	GLN	-	expression tag	UNP B8XCNO
D	656	PHE	-	expression tag	UNP B8XCNO
D	657	GLU	-	expression tag	UNP B8XCNO
D	658	LYS	-	expression tag	UNP B8XCNO
D	659	GLY	-	expression tag	UNP B8XCNO
D	660	GLY	-	expression tag	UNP B8XCNO
D	661	GLY	-	expression tag	UNP B8XCNO
D	662	SER	-	expression tag	UNP B8XCNO
D	663	GLY	-	expression tag	UNP B8XCNO

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Chain	Residue	Modelled	Actual	Comment	Reference
D	664	GLY	-	expression tag	UNP B8XCNO
D	665	GLY	-	expression tag	UNP B8XCNO
D	666	SER	-	expression tag	UNP B8XCNO
D	667	GLY	-	expression tag	UNP B8XCNO
D	668	GLY	-	expression tag	UNP B8XCNO
D	669	GLY	-	expression tag	UNP B8XCNO
D	670	SER	-	expression tag	UNP B8XCNO
D	671	TRP	-	expression tag	UNP B8XCNO
D	672	SER	-	expression tag	UNP B8XCNO
D	673	HIS	-	expression tag	UNP B8XCNO
D	674	PRO	-	expression tag	UNP B8XCNO
D	675	GLN	-	expression tag	UNP B8XCNO
D	676	PHE	-	expression tag	UNP B8XCNO
D	677	GLU	-	expression tag	UNP B8XCNO
D	678	LYS	-	expression tag	UNP B8XCNO
E	641	VAL	-	expression tag	UNP B8XCNO
E	642	GLU	-	expression tag	UNP B8XCNO
E	643	VAL	-	expression tag	UNP B8XCNO
E	644	ASP	-	expression tag	UNP B8XCNO
E	645	ASP	-	expression tag	UNP B8XCNO
E	646	ASP	-	expression tag	UNP B8XCNO
E	647	ASP	-	expression tag	UNP B8XCNO
E	648	LYS	-	expression tag	UNP B8XCNO
E	649	ALA	-	expression tag	UNP B8XCNO
E	650	GLY	-	expression tag	UNP B8XCNO
E	651	TRP	-	expression tag	UNP B8XCNO
E	652	SER	-	expression tag	UNP B8XCNO
E	653	HIS	-	expression tag	UNP B8XCNO
E	654	PRO	-	expression tag	UNP B8XCNO
E	655	GLN	-	expression tag	UNP B8XCNO
E	656	PHE	-	expression tag	UNP B8XCNO
E	657	GLU	-	expression tag	UNP B8XCNO
E	658	LYS	-	expression tag	UNP B8XCNO
E	659	GLY	-	expression tag	UNP B8XCNO
E	660	GLY	-	expression tag	UNP B8XCNO
E	661	GLY	-	expression tag	UNP B8XCNO
E	662	SER	-	expression tag	UNP B8XCNO
E	663	GLY	-	expression tag	UNP B8XCNO
E	664	GLY	-	expression tag	UNP B8XCNO
E	665	GLY	-	expression tag	UNP B8XCNO
E	666	SER	-	expression tag	UNP B8XCNO
E	667	GLY	-	expression tag	UNP B8XCNO

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Chain	Residue	Modelled	Actual	Comment	Reference
E	668	GLY	-	expression tag	UNP B8XCNO
E	669	GLY	-	expression tag	UNP B8XCNO
E	670	SER	-	expression tag	UNP B8XCNO
E	671	TRP	-	expression tag	UNP B8XCNO
E	672	SER	-	expression tag	UNP B8XCNO
E	673	HIS	-	expression tag	UNP B8XCNO
E	674	PRO	-	expression tag	UNP B8XCNO
E	675	GLN	-	expression tag	UNP B8XCNO
E	676	PHE	-	expression tag	UNP B8XCNO
E	677	GLU	-	expression tag	UNP B8XCNO
E	678	LYS	-	expression tag	UNP B8XCNO
F	641	VAL	-	expression tag	UNP B8XCNO
F	642	GLU	-	expression tag	UNP B8XCNO
F	643	VAL	-	expression tag	UNP B8XCNO
F	644	ASP	-	expression tag	UNP B8XCNO
F	645	ASP	-	expression tag	UNP B8XCNO
F	646	ASP	-	expression tag	UNP B8XCNO
F	647	ASP	-	expression tag	UNP B8XCNO
F	648	LYS	-	expression tag	UNP B8XCNO
F	649	ALA	-	expression tag	UNP B8XCNO
F	650	GLY	-	expression tag	UNP B8XCNO
F	651	TRP	-	expression tag	UNP B8XCNO
F	652	SER	-	expression tag	UNP B8XCNO
F	653	HIS	-	expression tag	UNP B8XCNO
F	654	PRO	-	expression tag	UNP B8XCNO
F	655	GLN	-	expression tag	UNP B8XCNO
F	656	PHE	-	expression tag	UNP B8XCNO
F	657	GLU	-	expression tag	UNP B8XCNO
F	658	LYS	-	expression tag	UNP B8XCNO
F	659	GLY	-	expression tag	UNP B8XCNO
F	660	GLY	-	expression tag	UNP B8XCNO
F	661	GLY	-	expression tag	UNP B8XCNO
F	662	SER	-	expression tag	UNP B8XCNO
F	663	GLY	-	expression tag	UNP B8XCNO
F	664	GLY	-	expression tag	UNP B8XCNO
F	665	GLY	-	expression tag	UNP B8XCNO
F	666	SER	-	expression tag	UNP B8XCNO
F	667	GLY	-	expression tag	UNP B8XCNO
F	668	GLY	-	expression tag	UNP B8XCNO
F	669	GLY	-	expression tag	UNP B8XCNO
F	670	SER	-	expression tag	UNP B8XCNO
F	671	TRP	-	expression tag	UNP B8XCNO

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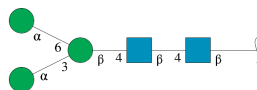
Chain	Residue	Modelled	Actual	Comment	Reference
F	672	SER	-	expression tag	UNP B8XCN0
F	673	HIS	-	expression tag	UNP B8XCN0
F	674	PRO	-	expression tag	UNP B8XCN0
F	675	GLN	-	expression tag	UNP B8XCN0
F	676	PHE	-	expression tag	UNP B8XCN0
F	677	GLU	-	expression tag	UNP B8XCN0
F	678	LYS	-	expression tag	UNP B8XCN0

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



Mol	Chain	Residues	Atoms				AltConf	Trace
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	N	2	Total	C	N	O	0	0
			28	16	2	10		
5	Q	2	Total	C	N	O	0	0
			28	16	2	10		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
6	O	5	Total	C	N	O	0	0
			61	34	2	25		
6	P	5	Total	C	N	O	0	0
			61	34	2	25		
6	R	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

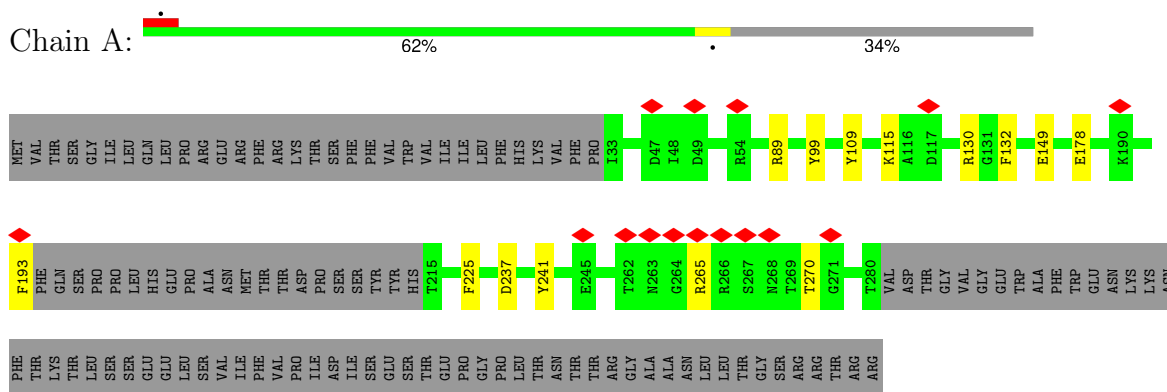


Mol	Chain	Residues	Atoms				AltConf
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	

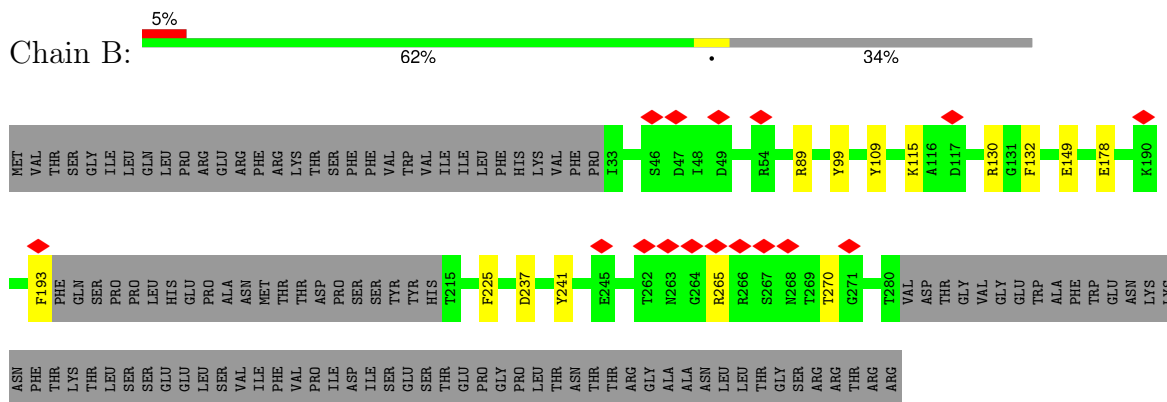
### 3 Residue-property plots [i](#)

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.

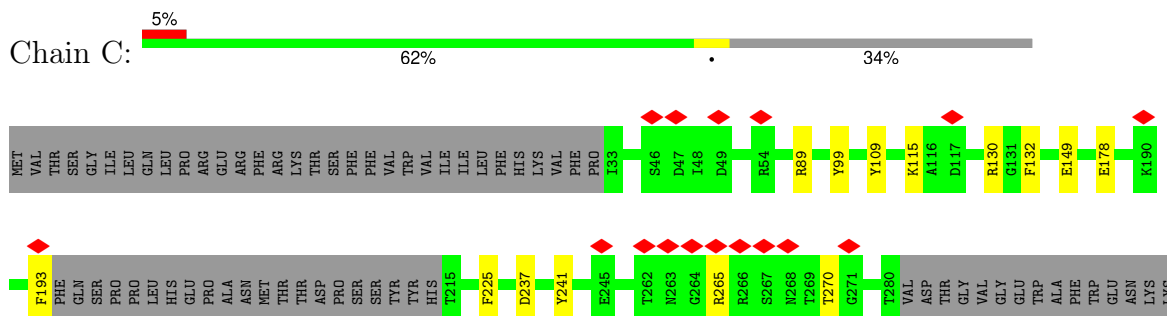
- Molecule 1: Spike glycoprotein 1



- Molecule 1: Spike glycoprotein 1



- Molecule 1: Spike glycoprotein 1



ASN PHE THR LYS THR LEU SER SER SER GLU LEU LEU SER SER VAL ILE PHE VAL VAL PRO ILE ASP ILE SER SER GLU THR THR THR ASN THR THR THR ARG GLY ALA ALA ASN LEU LEU THR THR GLY SER ARG THR ARG ARG

• Molecule 2: BDBV-43 Fab heavy chain

Chain G: 

MET GLU LEU GLY LEU ARG TRP PHE VAL LEU VAL ALA ILE LEU GLY VAL GLN CYS Q1 L4 E10 V11 R23 A24 S25 G26 D27 S28 R38 Q39 A40 Q43 G44 F45 A60 D72 E85 D86 D101 S102 W103 S112 SER ALA THR LYS GLY PRO

SER VAL PHE PRO LEU ALA PRO SER SER LEU LEU THR SER THR GLY THR GLY THR VAL ALA ALA LEU CYS GLY CYS VAL LYS ASP TYR PHE PRO VAL ASP PRO VAL ARG VAL THR VAL TRP TRP ASN SER GLY ALA LEU THR SER GLY VAL HIS THR PHE PRO VAL LEU GLN SER SER GLY TYR SER LEU SER

SER VAL VAL THR VAL PRO SER SER SER LEU LEU THR GLN THR TYR THR GLY THR ILE CYS VAL ASN VAL ASN HIS LYS PRO LYS ASN THR LYS VAL ASP PRO LYS ARG VAL GLU PRO LYS CYS SER ASP

• Molecule 2: BDBV-43 Fab heavy chain

Chain H: 

MET GLU LEU GLY LEU ARG TRP PHE VAL LEU VAL ALA ILE LEU GLU VAL GLN CYS Q1 L4 V11 R23 A24 S25 G26 D27 S28 R38 Q39 A40 Q43 G44 F45 A60 D72 L82 E85 D86 D101 S102 W103 S112 SER ALA THR LYS GLY

PRO SER VAL PHE PRO LEU ALA PRO SER SER LEU LEU THR SER THR GLY THR ILE CYS ALA LEU VAL ASN HIS LYS PRO LYS ASP TYR PHE PRO VAL ASP PRO VAL ARG VAL THR VAL TRP TRP ASN SER GLY ALA LEU THR SER GLY VAL HIS THR PHE PRO VAL LEU GLN SER SER GLY TYR SER LEU

SER SER VAL THR VAL PRO SER SER SER LEU LEU THR GLN THR TYR THR ILE CYS ALA VAL ASN VAL ASN HIS LYS PRO LYS ASP TYR PHE PRO VAL ASP PRO VAL ARG VAL THR VAL TRP TRP ASN SER GLY ALA LEU THR SER GLY VAL HIS THR PHE PRO VAL LEU GLN SER SER GLY TYR SER LEU

• Molecule 2: BDBV-43 Fab heavy chain


Chain I: 

MET GLU LEU GLY LEU ARG TRP PHE VAL LEU VAL ALA ILE LEU GLU VAL GLN CYS Q1 L4 V11 R23 A24 S25 G26 D27 S28 R38 Q39 A40 Q43 G44 F45 A60 D72 E85 D86 Y91 D101 S102 W103 S112 SER ALA THR LYS GLY

PRO SER VAL PHE PRO LEU ALA PRO SER SER LEU LEU THR SER THR GLY THR ILE CYS ALA LEU VAL ASN HIS LYS PRO LYS ASP TYR PHE PRO VAL ASP PRO VAL ARG VAL THR VAL TRP TRP ASN SER GLY ALA LEU THR SER GLY VAL HIS THR PHE PRO VAL LEU GLN SER SER GLY TYR SER LEU

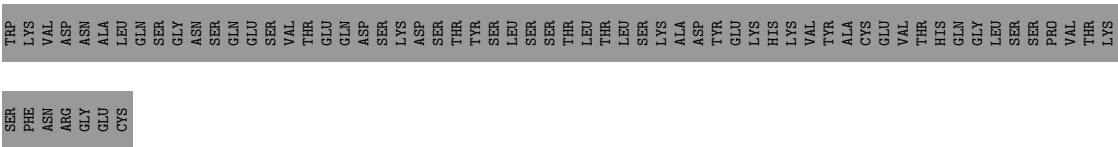
SER SER VAL THR VAL PRO SER SER SER LEU LEU THR GLN THR TYR THR ILE CYS ALA VAL ASN VAL ASN HIS LYS PRO LYS ASP TYR PHE PRO VAL ASP PRO VAL ARG VAL THR VAL TRP TRP ASN SER GLY ALA LEU THR SER GLY VAL HIS THR PHE PRO VAL LEU GLN SER SER GLY TYR SER LEU

• Molecule 3: BDBV-43 Fab light chain

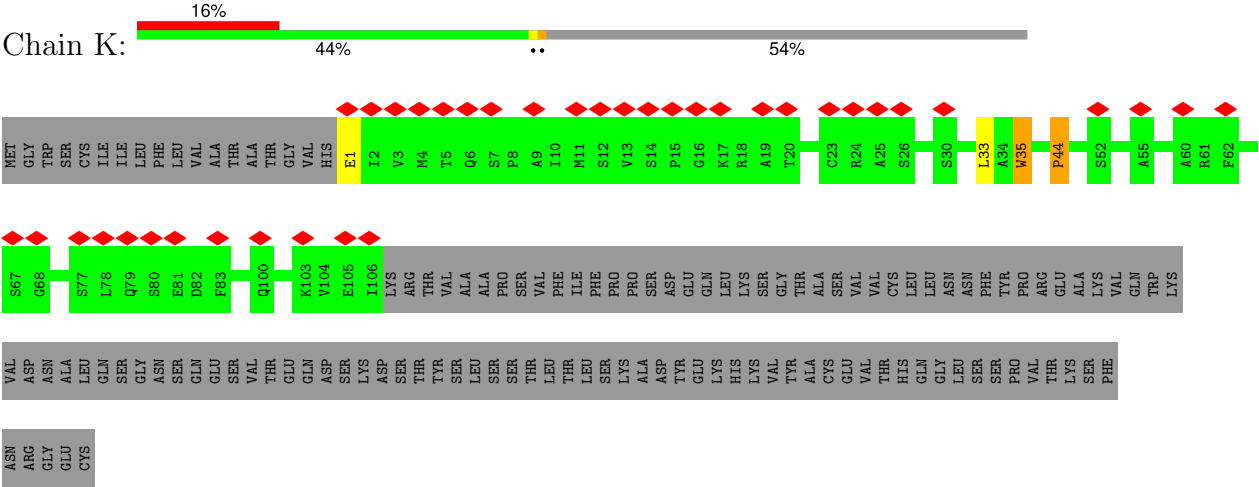
Chain J: 

MET GLY TRP SER CYS ILE ILE LEU PHE LEU VAL ALA THR ALA THR VAL HIS E1 I2 V3 M4 T5 Q6 S7 P8 A9 I10 M11 S12 V13 S14 P15 G16 K17 R18 A19 T20 C23 R24 A25 S26 S30 L33 A34 W35 P44 S52 A55 I58 P59 A60 R61

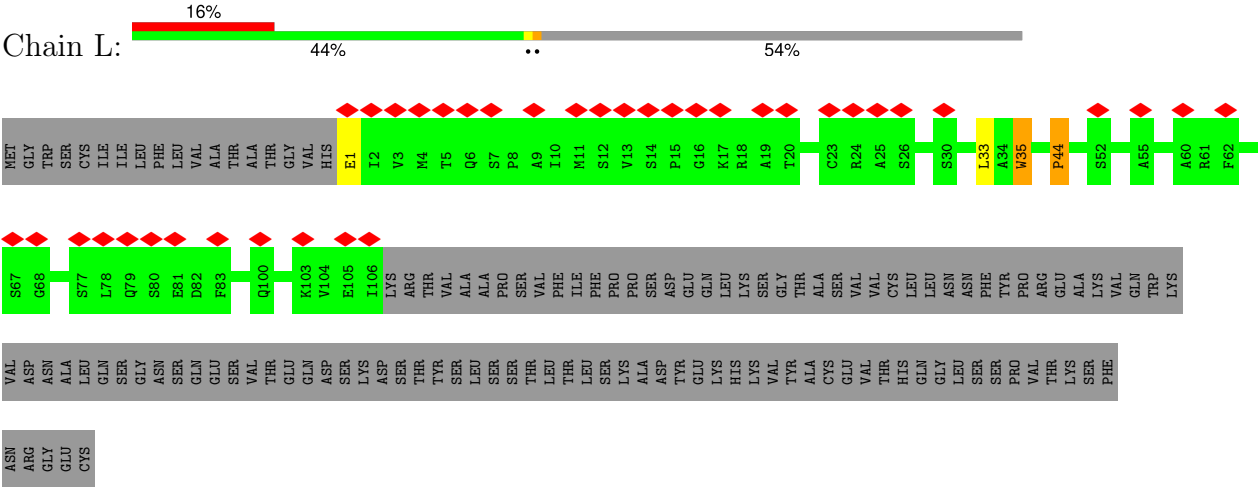
F62 S67 G68 S77 L78 Q79 S80 E81 D82 F83 Q100 K103 V104 E105 I106 LYS ARG THR VAL ALA ALA PRO PRO SER VAL PHE ILE PHE PRO PRO ASP ASP GLU GLN LEU LYS SER GLY THR ALA SER VAL CYS LEU LEU ASN ASN PHE TYR PRO ARG GLU ALA LYS VAL GLN



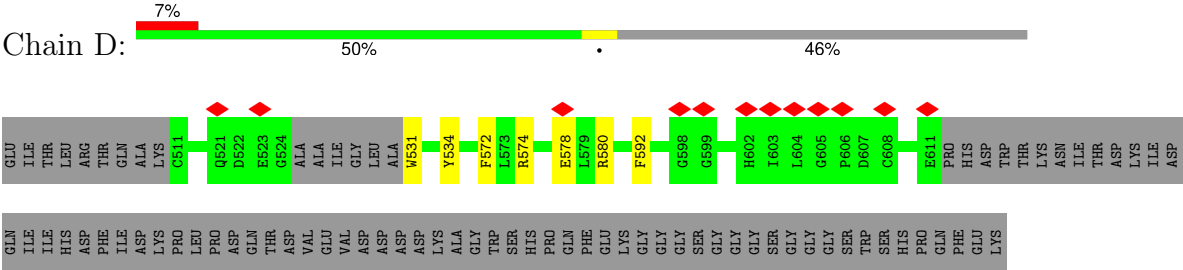
• Molecule 3: BDBV-43 Fab light chain



• Molecule 3: BDBV-43 Fab light chain

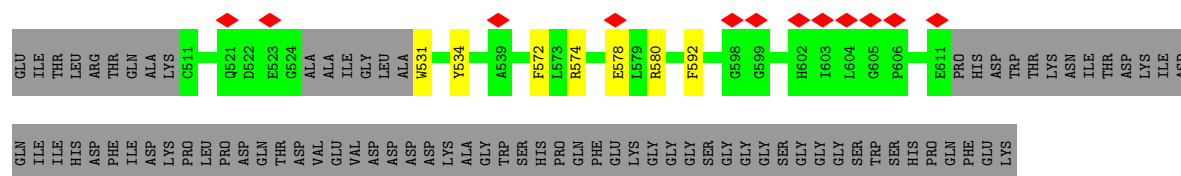


• Molecule 4: Envelope glycoprotein 2

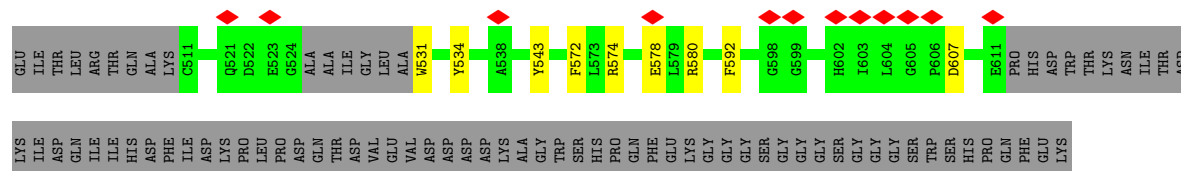


• Molecule 4: Envelope glycoprotein 2





• Molecule 4: Envelope glycoprotein 2



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



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



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



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



- Molecule 6:  $\alpha$ -D-mannopyranose-(1-3)-[ $\alpha$ -D-mannopyranose-(1-6)] $\beta$ -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- $\beta$ -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- $\beta$ -D-glucopyranose



- Molecule 6:  $\alpha$ -D-mannopyranose-(1-3)-[ $\alpha$ -D-mannopyranose-(1-6)] $\beta$ -D-mannopyranose-(1-4)-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	383783	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 ( $e^-/\text{\AA}^2$ )	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.150	Depositor
Minimum map value	-1.150	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.072	Depositor
Recommended contour level	0.41	Depositor
Map size (Å)	296.63998, 296.63998, 296.63998	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	1.13	6/1828 (0.3%)	0.96	6/2488 (0.2%)
1	B	1.12	6/1828 (0.3%)	0.96	6/2488 (0.2%)
1	C	1.13	6/1828 (0.3%)	0.95	6/2488 (0.2%)
2	G	1.15	5/951 (0.5%)	0.95	1/1288 (0.1%)
2	H	1.15	5/951 (0.5%)	0.95	1/1288 (0.1%)
2	I	1.15	5/951 (0.5%)	0.95	1/1288 (0.1%)
3	J	1.11	4/834 (0.5%)	0.95	0/1131
3	K	1.11	4/834 (0.5%)	0.95	0/1131
3	L	1.11	4/834 (0.5%)	0.95	0/1131
4	D	1.20	6/775 (0.8%)	1.00	4/1053 (0.4%)
4	E	1.20	6/775 (0.8%)	1.00	4/1053 (0.4%)
4	F	1.20	6/775 (0.8%)	1.00	5/1053 (0.5%)
All	All	1.14	63/13164 (0.5%)	0.96	34/17880 (0.2%)

All (63) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	578	GLU	CG-CD	7.70	1.63	1.51
4	E	578	GLU	CG-CD	7.64	1.63	1.51
4	D	578	GLU	CG-CD	7.60	1.63	1.51
1	B	193	PHE	CG-CD2	7.34	1.49	1.38
1	A	193	PHE	CG-CD2	7.23	1.49	1.38
1	C	193	PHE	CG-CD2	7.22	1.49	1.38
1	B	132	PHE	CB-CG	-6.67	1.40	1.51
1	C	132	PHE	CB-CG	-6.66	1.40	1.51
1	A	132	PHE	CB-CG	-6.62	1.40	1.51
2	H	1	GLN	CG-CD	6.35	1.65	1.51
2	I	1	GLN	CG-CD	6.33	1.65	1.51
2	G	1	GLN	CG-CD	6.30	1.65	1.51
3	L	35	TRP	CZ3-CH2	-6.27	1.30	1.40
3	J	35	TRP	CZ3-CH2	-6.26	1.30	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	193	PHE	CG-CD1	6.23	1.48	1.38
1	C	193	PHE	CG-CD1	6.20	1.48	1.38
4	F	572	PHE	CB-CG	-6.19	1.40	1.51
4	D	572	PHE	CB-CG	-6.18	1.40	1.51
4	E	572	PHE	CB-CG	-6.17	1.40	1.51
3	K	35	TRP	CZ3-CH2	-6.17	1.30	1.40
1	A	193	PHE	CG-CD1	6.16	1.48	1.38
2	H	103	TRP	CB-CG	-6.04	1.39	1.50
2	G	103	TRP	CB-CG	-5.96	1.39	1.50
2	I	103	TRP	CB-CG	-5.96	1.39	1.50
3	K	35	TRP	CD2-CE2	-5.95	1.34	1.41
3	J	35	TRP	CD2-CE2	-5.91	1.34	1.41
1	C	178	GLU	CD-OE2	-5.91	1.19	1.25
3	L	35	TRP	CD2-CE2	-5.89	1.34	1.41
3	J	1	GLU	CB-CG	5.86	1.63	1.52
3	K	1	GLU	CB-CG	5.85	1.63	1.52
1	A	241	TYR	CB-CG	-5.83	1.43	1.51
3	L	1	GLU	CB-CG	5.82	1.63	1.52
1	C	241	TYR	CB-CG	-5.75	1.43	1.51
1	B	241	TYR	CB-CG	-5.74	1.43	1.51
1	B	178	GLU	CD-OE2	-5.73	1.19	1.25
2	I	103	TRP	NE1-CE2	-5.65	1.30	1.37
4	D	531	TRP	CZ2-CH2	5.62	1.48	1.37
1	A	178	GLU	CD-OE2	-5.60	1.19	1.25
4	E	531	TRP	CZ2-CH2	5.58	1.48	1.37
2	G	103	TRP	NE1-CE2	-5.58	1.30	1.37
4	F	531	TRP	CZ2-CH2	5.58	1.48	1.37
2	H	103	TRP	NE1-CE2	-5.58	1.30	1.37
4	E	592	PHE	CB-CG	-5.56	1.42	1.51
4	D	531	TRP	CD2-CE3	5.54	1.48	1.40
4	D	592	PHE	CB-CG	-5.51	1.42	1.51
4	E	531	TRP	CD2-CE3	5.50	1.48	1.40
4	F	531	TRP	CD2-CE3	5.48	1.48	1.40
3	K	44	PRO	N-CD	-5.47	1.40	1.47
3	J	44	PRO	N-CD	-5.46	1.40	1.47
3	L	44	PRO	N-CD	-5.46	1.40	1.47
4	F	592	PHE	CB-CG	-5.43	1.42	1.51
1	A	193	PHE	CE2-CZ	5.13	1.47	1.37
4	F	531	TRP	CE3-CZ3	5.13	1.47	1.38
2	H	103	TRP	CE2-CZ2	-5.12	1.31	1.39
1	C	193	PHE	CE2-CZ	5.11	1.47	1.37
2	G	103	TRP	CE2-CZ2	-5.10	1.31	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	531	TRP	CE3-CZ3	5.10	1.47	1.38
1	B	193	PHE	CE2-CZ	5.09	1.47	1.37
2	I	103	TRP	CE2-CZ2	-5.06	1.31	1.39
4	E	531	TRP	CE3-CZ3	5.03	1.47	1.38
2	G	10	GLU	CD-OE2	-5.03	1.20	1.25
2	H	103	TRP	CD2-CE2	-5.01	1.35	1.41
2	I	91	TYR	CB-CG	-5.01	1.44	1.51

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	534	TYR	CB-CG-CD1	-8.93	115.64	121.00
4	F	534	TYR	CB-CG-CD1	-8.88	115.67	121.00
4	D	534	TYR	CB-CG-CD1	-8.79	115.73	121.00
1	B	89	ARG	NE-CZ-NH2	-7.75	116.42	120.30
1	A	89	ARG	NE-CZ-NH2	-7.73	116.44	120.30
1	C	89	ARG	NE-CZ-NH2	-7.68	116.46	120.30
4	D	574	ARG	NE-CZ-NH2	-7.25	116.68	120.30
4	F	574	ARG	NE-CZ-NH2	-7.00	116.80	120.30
4	E	574	ARG	NE-CZ-NH2	-6.91	116.84	120.30
4	F	534	TYR	CA-CB-CG	6.32	125.40	113.40
1	B	109	TYR	CB-CG-CD1	-6.30	117.22	121.00
4	D	534	TYR	CA-CB-CG	6.27	125.31	113.40
4	E	534	TYR	CA-CB-CG	6.26	125.30	113.40
1	A	109	TYR	CB-CG-CD1	-6.25	117.25	121.00
1	C	109	TYR	CB-CG-CD1	-6.20	117.28	121.00
1	A	130	ARG	NE-CZ-NH2	-6.15	117.23	120.30
1	C	130	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	B	130	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	C	265	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	B	265	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	265	ARG	NE-CZ-NH1	5.58	123.09	120.30
2	G	38	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	C	99	TYR	CB-CG-CD1	-5.49	117.71	121.00
1	B	99	TYR	CB-CG-CD1	-5.49	117.71	121.00
2	I	38	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	A	99	TYR	CB-CG-CD1	-5.43	117.75	121.00
2	H	38	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	B	237	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	237	ASP	CB-CG-OD2	5.33	123.10	118.30
4	D	580	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	C	237	ASP	CB-CG-OD2	5.24	123.02	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	580	ARG	NE-CZ-NH1	5.09	122.85	120.30
4	F	543	TYR	CB-CG-CD1	-5.01	117.99	121.00
4	E	580	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	1785	0	1747	1	0
1	B	1785	0	1747	1	0
1	C	1785	0	1747	1	0
2	G	931	0	913	2	0
2	H	931	0	913	2	0
2	I	931	0	913	2	0
3	J	815	0	804	3	0
3	K	815	0	804	2	0
3	L	815	0	804	2	0
4	D	757	0	723	0	0
4	E	757	0	723	0	0
4	F	757	0	723	1	0
5	M	28	0	25	0	0
5	N	28	0	25	0	0
5	Q	28	0	25	0	0
6	O	61	0	52	0	0
6	P	61	0	52	0	0
6	R	61	0	52	0	0
7	A	14	0	13	0	0
7	B	14	0	13	0	0
7	C	14	0	13	0	0
All	All	13173	0	12831	14	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 1.

All (14) 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)
2:I:103:TRP:CD1	3:L:44:PRO:HD2	2.42	0.55
2:G:103:TRP:CD1	3:J:44:PRO:HD2	2.42	0.55
2:H:103:TRP:CD1	3:K:44:PRO:HD2	2.42	0.55
2:G:27:ASP:OD1	2:G:27:ASP:N	2.45	0.48
2:I:27:ASP:OD1	2:I:27:ASP:N	2.45	0.44
2:H:27:ASP:OD1	2:H:27:ASP:N	2.45	0.43
3:L:33:LEU:HB2	3:L:35:TRP:CH2	2.54	0.43
3:K:33:LEU:HB2	3:K:35:TRP:CH2	2.54	0.42
3:J:58:ILE:HA	3:J:59:PRO:HD3	1.94	0.42
3:J:33:LEU:HB2	3:J:35:TRP:CH2	2.54	0.42
1:B:115:LYS:NZ	1:B:149:GLU:OE2	2.54	0.41
1:C:115:LYS:NZ	1:C:149:GLU:OE2	2.53	0.41
1:A:115:LYS:NZ	1:A:149:GLU:OE2	2.53	0.41
4:F:607:ASP:N	4:F:607:ASP:OD1	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	223/343 (65%)	211 (95%)	10 (4%)	2 (1%)	14	50
1	B	223/343 (65%)	211 (95%)	10 (4%)	2 (1%)	14	50
1	C	223/343 (65%)	211 (95%)	10 (4%)	2 (1%)	14	50
2	G	120/246 (49%)	118 (98%)	2 (2%)	0	100	100
2	H	120/246 (49%)	118 (98%)	2 (2%)	0	100	100
2	I	120/246 (49%)	118 (98%)	2 (2%)	0	100	100
3	J	104/232 (45%)	101 (97%)	3 (3%)	0	100	100
3	K	104/232 (45%)	101 (97%)	3 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	104/232 (45%)	101 (97%)	3 (3%)	0	100	100
4	D	91/177 (51%)	89 (98%)	2 (2%)	0	100	100
4	E	91/177 (51%)	89 (98%)	2 (2%)	0	100	100
4	F	91/177 (51%)	89 (98%)	2 (2%)	0	100	100
All	All	1614/2994 (54%)	1557 (96%)	51 (3%)	6 (0%)	32	67

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	THR
1	B	270	THR
1	C	270	THR
1	A	225	PHE
1	B	225	PHE
1	C	225	PHE

### 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	197/303 (65%)	197 (100%)	0	100	100
1	B	197/303 (65%)	197 (100%)	0	100	100
1	C	197/303 (65%)	197 (100%)	0	100	100
2	G	101/208 (49%)	101 (100%)	0	100	100
2	H	101/208 (49%)	101 (100%)	0	100	100
2	I	101/208 (49%)	101 (100%)	0	100	100
3	J	89/200 (44%)	89 (100%)	0	100	100
3	K	89/200 (44%)	89 (100%)	0	100	100
3	L	89/200 (44%)	89 (100%)	0	100	100
4	D	80/146 (55%)	80 (100%)	0	100	100
4	E	80/146 (55%)	80 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	F	80/146 (55%)	80 (100%)	0	100	100
All	All	1401/2571 (54%)	1401 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

21 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	M	1	5,1	14,14,15	2.15	7 (50%)	17,19,21	1.36	4 (23%)
5	NAG	M	2	5	14,14,15	2.14	6 (42%)	17,19,21	1.08	2 (11%)
5	NAG	N	1	5,1	14,14,15	2.17	7 (50%)	17,19,21	1.36	4 (23%)
5	NAG	N	2	5	14,14,15	2.17	5 (35%)	17,19,21	1.05	2 (11%)
6	NAG	O	1	4,6	14,14,15	2.17	7 (50%)	17,19,21	1.04	1 (5%)
6	NAG	O	2	6	14,14,15	1.89	4 (28%)	17,19,21	0.94	1 (5%)
6	BMA	O	3	6	11,11,12	1.41	3 (27%)	15,15,17	0.59	0
6	MAN	O	4	6	11,11,12	1.87	4 (36%)	15,15,17	0.64	0
6	MAN	O	5	6	11,11,12	1.83	4 (36%)	15,15,17	0.64	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	NAG	P	1	4,6	14,14,15	2.17	6 (42%)	17,19,21	1.06	1 (5%)
6	NAG	P	2	6	14,14,15	1.90	4 (28%)	17,19,21	0.93	1 (5%)
6	BMA	P	3	6	11,11,12	1.41	3 (27%)	15,15,17	0.58	0
6	MAN	P	4	6	11,11,12	1.87	5 (45%)	15,15,17	0.63	0
6	MAN	P	5	6	11,11,12	1.86	5 (45%)	15,15,17	0.65	0
5	NAG	Q	1	5,1	14,14,15	2.16	7 (50%)	17,19,21	1.35	4 (23%)
5	NAG	Q	2	5	14,14,15	2.15	5 (35%)	17,19,21	1.06	2 (11%)
6	NAG	R	1	4,6	14,14,15	2.15	7 (50%)	17,19,21	1.05	1 (5%)
6	NAG	R	2	6	14,14,15	1.91	4 (28%)	17,19,21	0.94	1 (5%)
6	BMA	R	3	6	11,11,12	1.41	3 (27%)	15,15,17	0.59	0
6	MAN	R	4	6	11,11,12	1.87	4 (36%)	15,15,17	0.64	0
6	MAN	R	5	6	11,11,12	1.86	5 (45%)	15,15,17	0.66	0

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	M	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	0/6/23/26	0/1/1/1
5	NAG	N	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	N	2	5	-	1/6/23/26	0/1/1/1
6	NAG	O	1	4,6	-	0/6/23/26	0/1/1/1
6	NAG	O	2	6	-	0/6/23/26	0/1/1/1
6	BMA	O	3	6	-	0/2/19/22	0/1/1/1
6	MAN	O	4	6	-	0/2/19/22	0/1/1/1
6	MAN	O	5	6	-	0/2/19/22	0/1/1/1
6	NAG	P	1	4,6	-	0/6/23/26	0/1/1/1
6	NAG	P	2	6	-	0/6/23/26	0/1/1/1
6	BMA	P	3	6	-	0/2/19/22	0/1/1/1
6	MAN	P	4	6	-	0/2/19/22	0/1/1/1
6	MAN	P	5	6	-	0/2/19/22	0/1/1/1
5	NAG	Q	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	0/6/23/26	0/1/1/1
6	NAG	R	1	4,6	-	0/6/23/26	0/1/1/1
6	NAG	R	2	6	-	0/6/23/26	0/1/1/1
6	BMA	R	3	6	-	0/2/19/22	0/1/1/1
6	MAN	R	4	6	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	MAN	R	5	6	-	0/2/19/22	0/1/1/1

All (105) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	O	1	NAG	C1-C2	5.60	1.60	1.52
6	P	1	NAG	C1-C2	5.60	1.60	1.52
5	N	2	NAG	C1-C2	5.59	1.60	1.52
5	Q	2	NAG	C1-C2	5.56	1.59	1.52
5	M	2	NAG	C1-C2	5.44	1.59	1.52
6	R	1	NAG	C1-C2	5.44	1.59	1.52
5	Q	1	NAG	C1-C2	5.29	1.59	1.52
5	N	1	NAG	C1-C2	5.28	1.59	1.52
5	M	1	NAG	C1-C2	5.24	1.59	1.52
6	R	2	NAG	C1-C2	4.42	1.58	1.52
6	O	2	NAG	C1-C2	4.36	1.58	1.52
6	P	2	NAG	C1-C2	4.31	1.58	1.52
6	P	5	MAN	O5-C5	3.24	1.49	1.43
6	O	5	MAN	O5-C5	3.19	1.49	1.43
6	P	4	MAN	O5-C5	3.18	1.49	1.43
6	R	5	MAN	O5-C5	3.18	1.49	1.43
6	O	4	MAN	O5-C5	3.14	1.49	1.43
6	R	4	MAN	O5-C5	3.10	1.49	1.43
5	N	2	NAG	O5-C5	2.98	1.49	1.43
5	M	2	NAG	O5-C5	2.97	1.49	1.43
5	Q	2	NAG	O5-C5	2.96	1.49	1.43
6	P	5	MAN	C2-C3	2.95	1.57	1.52
6	R	1	NAG	O5-C5	2.95	1.49	1.43
6	R	5	MAN	C2-C3	2.95	1.57	1.52
6	O	1	NAG	O5-C5	2.94	1.49	1.43
6	P	2	NAG	O5-C5	2.93	1.49	1.43
6	P	1	NAG	O5-C5	2.90	1.49	1.43
6	R	4	MAN	C2-C3	2.89	1.56	1.52
6	O	4	MAN	C2-C3	2.89	1.56	1.52
6	P	4	MAN	C2-C3	2.88	1.56	1.52
6	O	2	NAG	O5-C5	2.87	1.49	1.43
5	Q	1	NAG	O5-C5	2.86	1.49	1.43
5	N	1	NAG	O5-C5	2.86	1.49	1.43
6	O	5	MAN	C2-C3	2.83	1.56	1.52
5	M	1	NAG	O5-C5	2.82	1.48	1.43
6	R	2	NAG	O5-C5	2.81	1.48	1.43
5	N	2	NAG	C2-N2	2.55	1.50	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	2	NAG	C2-N2	2.55	1.50	1.46
5	N	2	NAG	C3-C2	2.54	1.57	1.52
6	R	5	MAN	C1-C2	2.51	1.58	1.52
5	M	2	NAG	C3-C2	2.51	1.57	1.52
6	P	4	MAN	C1-C2	2.50	1.58	1.52
6	R	4	MAN	C1-C2	2.50	1.58	1.52
6	O	4	MAN	C1-C2	2.50	1.58	1.52
5	Q	2	NAG	C3-C2	2.49	1.57	1.52
6	P	2	NAG	C3-C2	2.48	1.57	1.52
6	R	2	NAG	C3-C2	2.47	1.57	1.52
5	Q	2	NAG	C2-N2	2.46	1.50	1.46
6	P	5	MAN	C1-C2	2.44	1.58	1.52
6	O	2	NAG	C3-C2	2.43	1.57	1.52
6	O	5	MAN	C1-C2	2.43	1.58	1.52
6	R	4	MAN	C4-C5	2.40	1.58	1.53
6	O	4	MAN	C4-C5	2.39	1.58	1.53
5	N	1	NAG	C3-C2	2.38	1.57	1.52
6	P	1	NAG	C3-C2	2.37	1.57	1.52
5	M	1	NAG	C3-C2	2.37	1.57	1.52
6	P	5	MAN	C4-C5	2.37	1.58	1.53
6	R	5	MAN	C4-C5	2.37	1.58	1.53
5	Q	1	NAG	C3-C2	2.36	1.57	1.52
5	N	1	NAG	O5-C1	2.34	1.47	1.43
6	O	5	MAN	C4-C5	2.34	1.58	1.53
6	P	3	BMA	C2-C3	2.32	1.56	1.52
6	R	3	BMA	O5-C5	2.32	1.48	1.43
5	N	1	NAG	C2-N2	2.32	1.50	1.46
6	P	4	MAN	C4-C5	2.31	1.58	1.53
5	M	1	NAG	O5-C1	2.31	1.47	1.43
6	P	3	BMA	O5-C5	2.30	1.47	1.43
6	R	1	NAG	C3-C2	2.29	1.57	1.52
5	Q	1	NAG	C2-N2	2.29	1.50	1.46
6	R	3	BMA	C2-C3	2.28	1.56	1.52
6	O	3	BMA	C2-C3	2.27	1.56	1.52
6	R	1	NAG	C4-C3	2.26	1.58	1.52
6	O	1	NAG	C4-C3	2.26	1.58	1.52
6	O	1	NAG	C3-C2	2.26	1.57	1.52
6	O	3	BMA	O5-C5	2.26	1.47	1.43
5	M	1	NAG	C2-N2	2.25	1.50	1.46
5	M	1	NAG	C4-C5	2.25	1.57	1.53
5	N	1	NAG	C4-C5	2.25	1.57	1.53
5	Q	1	NAG	O5-C1	2.23	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	P	1	NAG	C4-C3	2.21	1.58	1.52
5	Q	1	NAG	C4-C5	2.21	1.57	1.53
6	O	1	NAG	C4-C5	2.20	1.57	1.53
6	R	1	NAG	O5-C1	2.18	1.47	1.43
6	O	3	BMA	C1-C2	2.18	1.57	1.52
6	R	2	NAG	C4-C5	2.17	1.57	1.53
6	R	3	BMA	C1-C2	2.17	1.57	1.52
6	P	1	NAG	O5-C1	2.17	1.47	1.43
5	N	2	NAG	C4-C5	2.15	1.57	1.53
5	Q	2	NAG	C4-C5	2.15	1.57	1.53
5	M	2	NAG	C4-C5	2.15	1.57	1.53
6	P	1	NAG	C4-C5	2.15	1.57	1.53
6	P	3	BMA	C1-C2	2.14	1.57	1.52
6	O	1	NAG	O5-C1	2.13	1.47	1.43
6	P	2	NAG	C4-C5	2.13	1.57	1.53
6	O	2	NAG	C4-C5	2.10	1.57	1.53
6	R	1	NAG	C4-C5	2.10	1.57	1.53
5	M	1	NAG	C4-C3	2.06	1.57	1.52
5	N	1	NAG	C4-C3	2.06	1.57	1.52
5	Q	1	NAG	C4-C3	2.05	1.57	1.52
6	R	5	MAN	C4-C3	2.05	1.57	1.52
6	P	5	MAN	C4-C3	2.04	1.57	1.52
6	R	1	NAG	C2-N2	2.03	1.49	1.46
6	P	4	MAN	C4-C3	2.03	1.57	1.52
5	M	2	NAG	C4-C3	2.02	1.57	1.52
6	O	1	NAG	C2-N2	2.01	1.49	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Q	1	NAG	C8-C7-N2	2.95	121.00	116.12
5	N	1	NAG	C8-C7-N2	2.94	121.00	116.12
5	M	1	NAG	C8-C7-N2	2.92	120.96	116.12
5	M	2	NAG	C8-C7-N2	2.83	120.82	116.12
5	N	2	NAG	C8-C7-N2	2.74	120.67	116.12
5	Q	2	NAG	C8-C7-N2	2.74	120.67	116.12
6	P	1	NAG	C8-C7-N2	2.50	120.26	116.12
6	R	1	NAG	C8-C7-N2	2.45	120.19	116.12
6	O	1	NAG	C8-C7-N2	2.42	120.12	116.12
5	Q	1	NAG	O7-C7-C8	-2.39	117.80	122.05
5	N	1	NAG	O7-C7-C8	-2.36	117.86	122.05
5	M	1	NAG	O7-C7-C8	-2.35	117.87	122.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Q	1	NAG	O4-C4-C3	-2.20	105.18	110.38
5	N	1	NAG	O4-C4-C3	-2.19	105.21	110.38
6	O	2	NAG	C8-C7-N2	2.19	119.74	116.12
5	M	2	NAG	O7-C7-C8	-2.18	118.17	122.05
5	M	1	NAG	O4-C4-C3	-2.18	105.23	110.38
5	Q	2	NAG	O7-C7-C8	-2.18	118.18	122.05
6	P	2	NAG	C8-C7-N2	2.17	119.72	116.12
6	R	2	NAG	C8-C7-N2	2.17	119.71	116.12
5	N	1	NAG	O5-C1-C2	-2.15	107.97	111.29
5	N	2	NAG	O7-C7-C8	-2.11	118.29	122.05
5	M	1	NAG	O5-C1-C2	-2.10	108.04	111.29
5	Q	1	NAG	O5-C1-C2	-2.05	108.11	111.29

There are no chirality outliers.

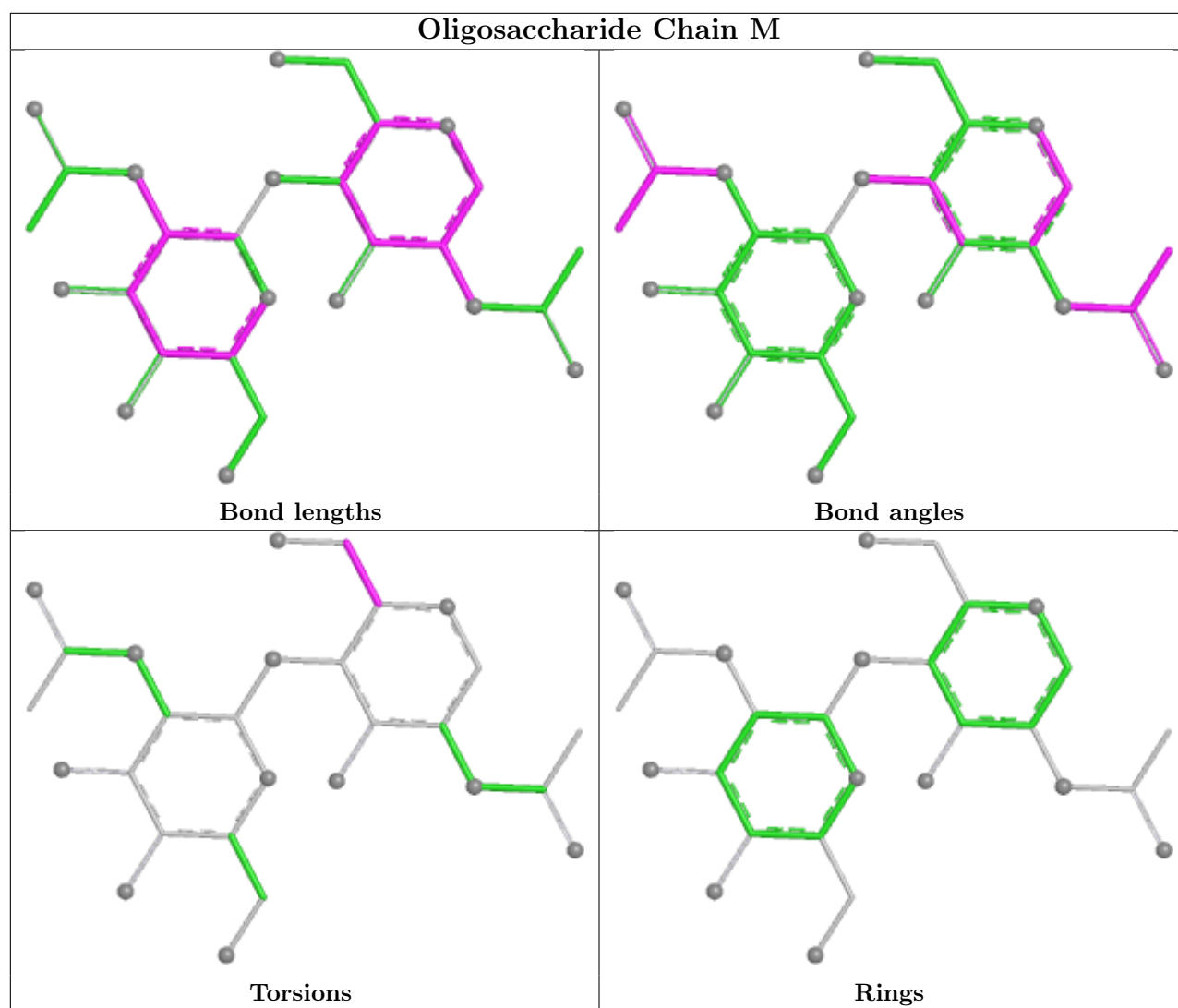
All (7) torsion outliers are listed below:

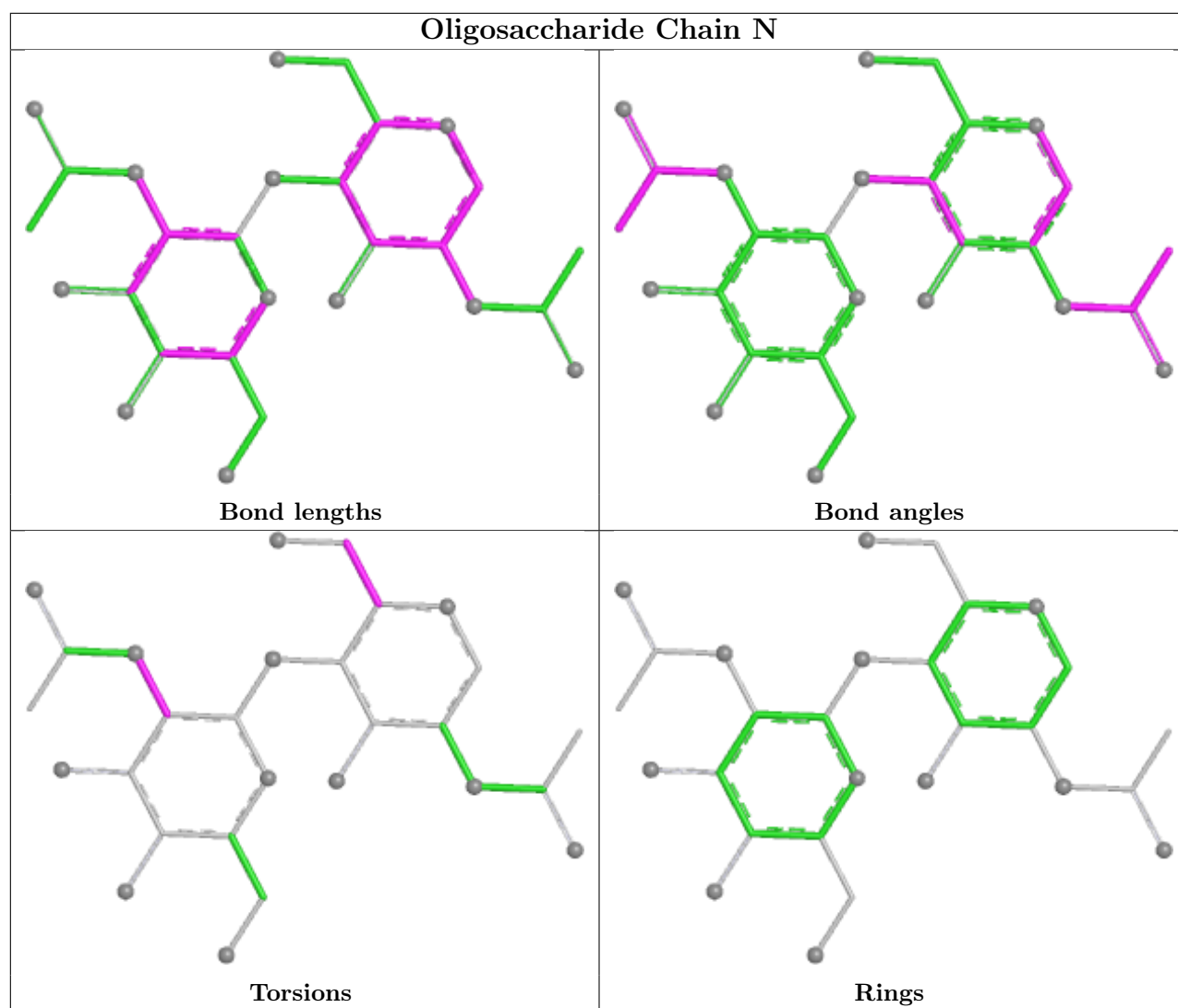
Mol	Chain	Res	Type	Atoms
5	N	1	NAG	C4-C5-C6-O6
5	Q	1	NAG	C4-C5-C6-O6
5	M	1	NAG	C4-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6
5	N	1	NAG	O5-C5-C6-O6
5	Q	1	NAG	O5-C5-C6-O6
5	N	2	NAG	C1-C2-N2-C7

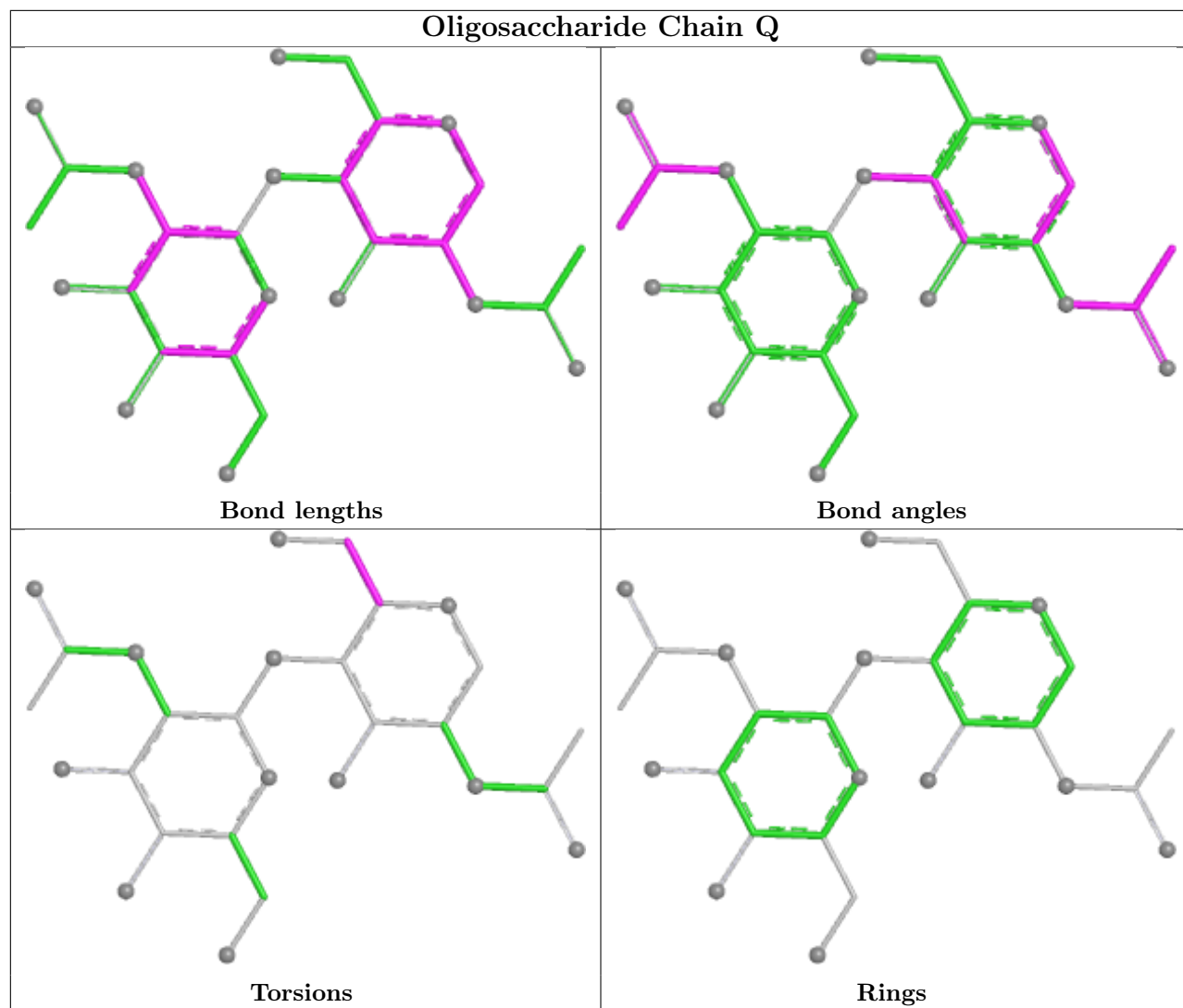
There are no ring outliers.

No monomer is involved in short contacts.

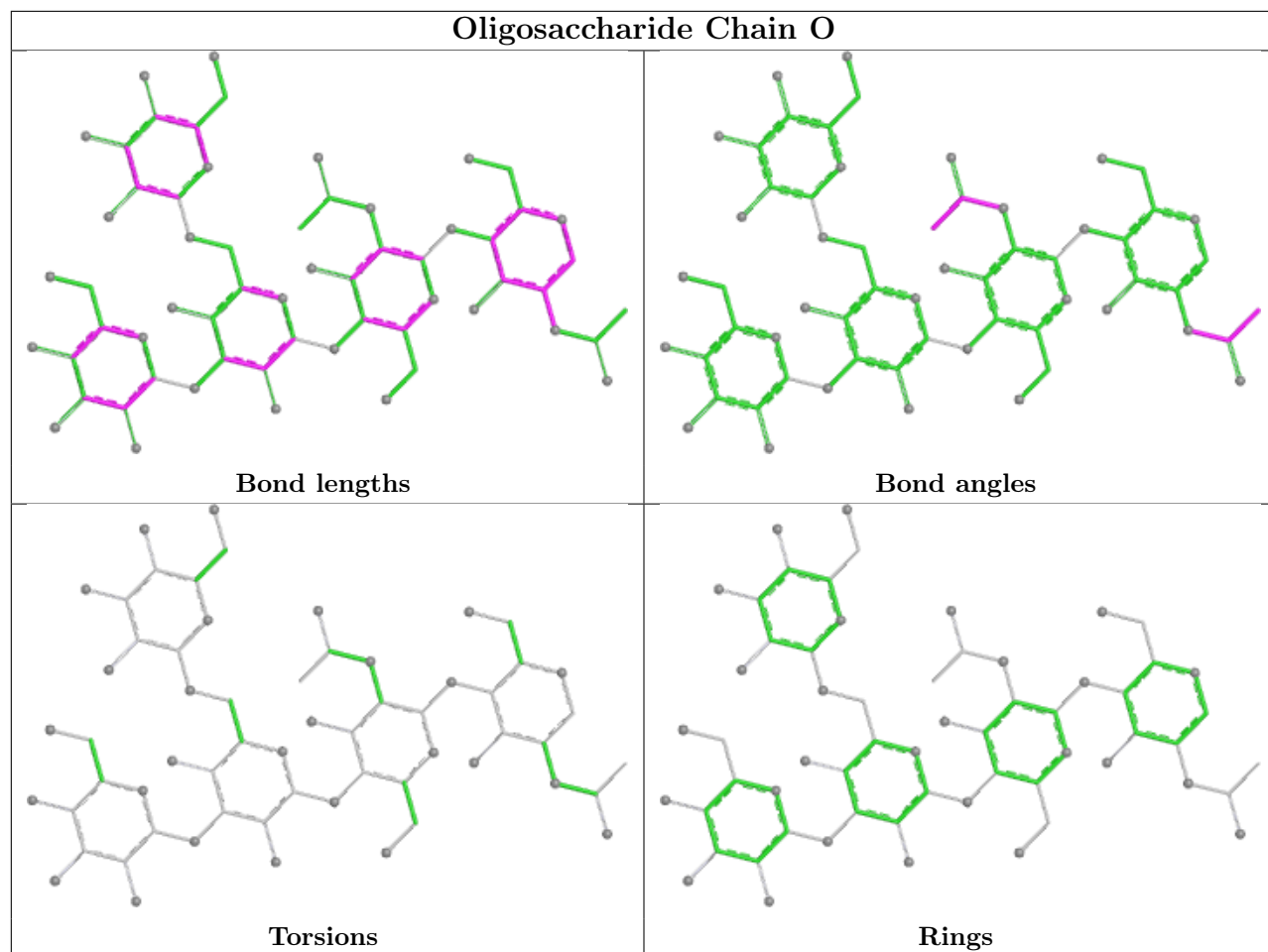
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

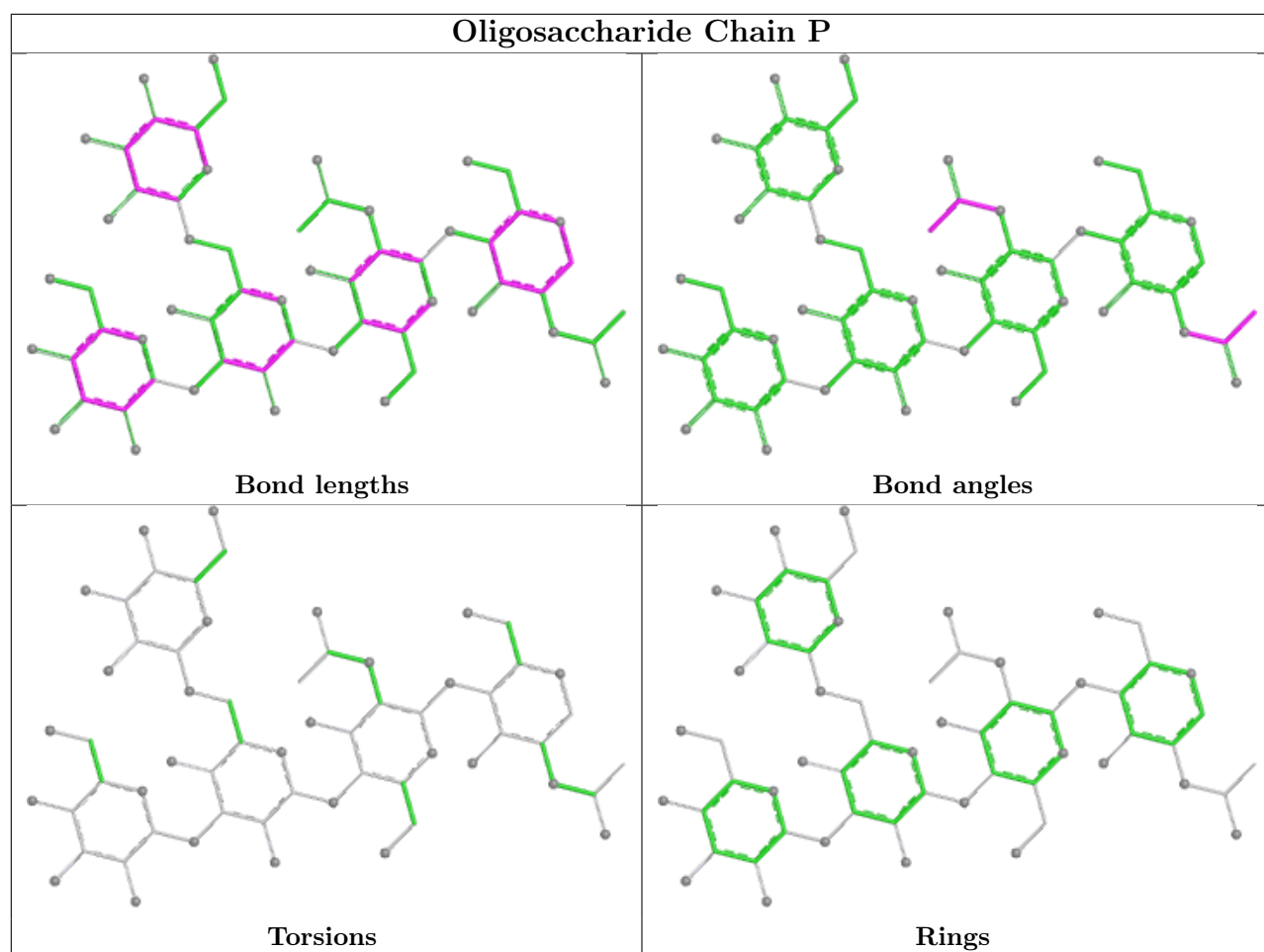


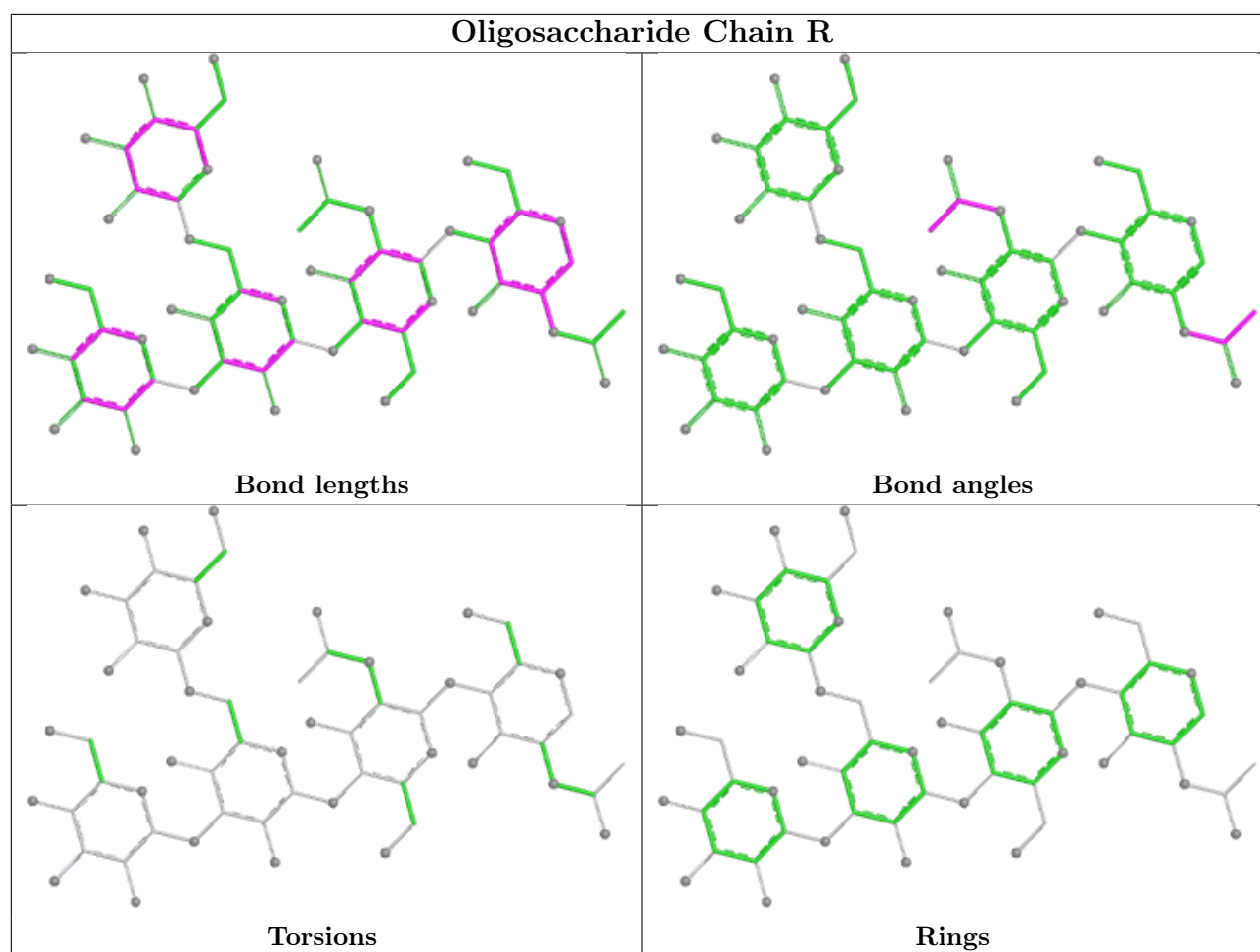












## 5.6 Ligand geometry [i](#)

3 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	$\# Z  > 2$
7	NAG	C	401	1	14,14,15	2.26	6 (42%)	17,19,21	1.03	1 (5%)
7	NAG	A	401	1	14,14,15	2.27	6 (42%)	17,19,21	1.02	2 (11%)
7	NAG	B	401	1	14,14,15	2.30	6 (42%)	17,19,21	1.02	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
7	NAG	C	401	1	-	0/6/23/26	0/1/1/1
7	NAG	A	401	1	-	0/6/23/26	0/1/1/1
7	NAG	B	401	1	-	0/6/23/26	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	401	NAG	C1-C2	5.76	1.60	1.52
7	A	401	NAG	C1-C2	5.70	1.60	1.52
7	C	401	NAG	C1-C2	5.63	1.60	1.52
7	B	401	NAG	O5-C5	3.51	1.50	1.43
7	A	401	NAG	O5-C5	3.51	1.50	1.43
7	C	401	NAG	O5-C5	3.46	1.50	1.43
7	B	401	NAG	C3-C2	2.58	1.57	1.52
7	B	401	NAG	O5-C1	2.56	1.48	1.43
7	C	401	NAG	C3-C2	2.53	1.57	1.52
7	A	401	NAG	C3-C2	2.52	1.57	1.52
7	C	401	NAG	O5-C1	2.51	1.47	1.43
7	A	401	NAG	O5-C1	2.47	1.47	1.43
7	A	401	NAG	C4-C5	2.33	1.58	1.53
7	B	401	NAG	C4-C5	2.30	1.57	1.53
7	C	401	NAG	C4-C5	2.30	1.57	1.53
7	C	401	NAG	C4-C3	2.24	1.58	1.52
7	B	401	NAG	C4-C3	2.19	1.58	1.52
7	A	401	NAG	C4-C3	2.16	1.58	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	401	NAG	C8-C7-N2	2.28	119.89	116.12
7	A	401	NAG	C8-C7-N2	2.22	119.80	116.12
7	B	401	NAG	C8-C7-N2	2.22	119.79	116.12
7	A	401	NAG	O7-C7-C8	-2.01	118.47	122.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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.

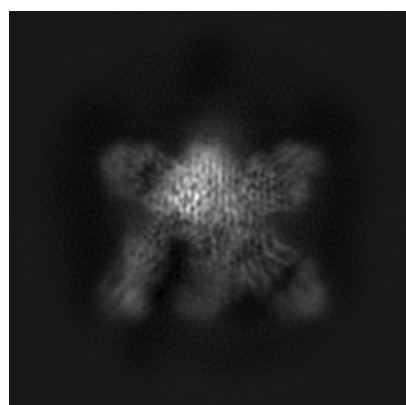
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22841. These allow visual inspection of the internal detail of the map and identification of artifacts.

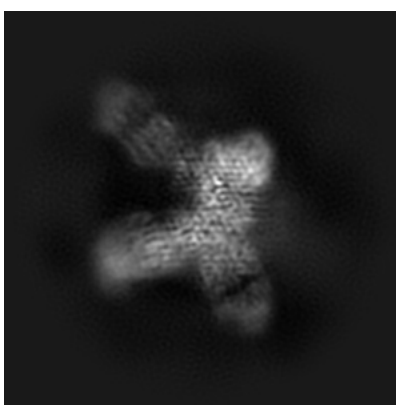
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

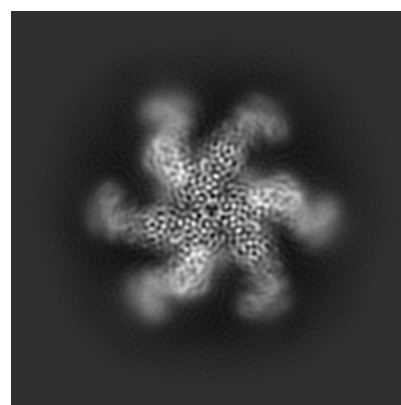
#### 6.1.1 Primary map



X



Y

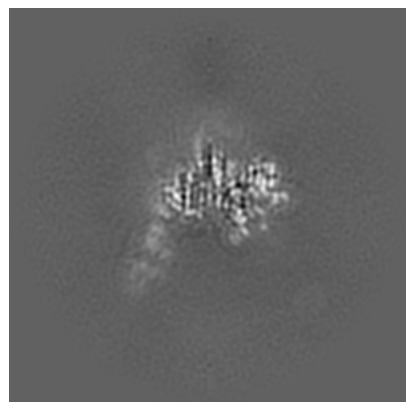


Z

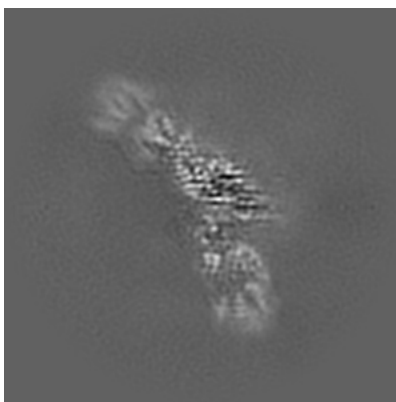
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

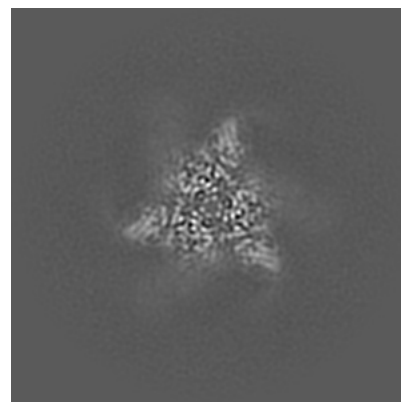
#### 6.2.1 Primary map



X Index: 144



Y Index: 144

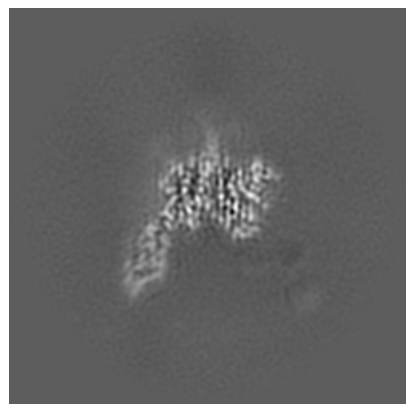


Z Index: 144

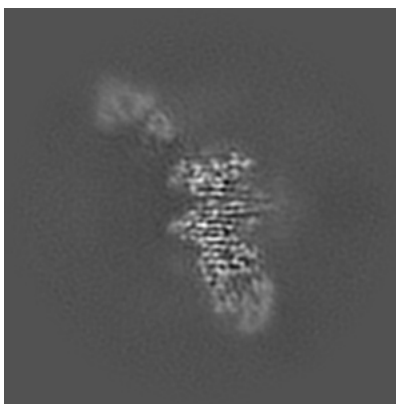
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

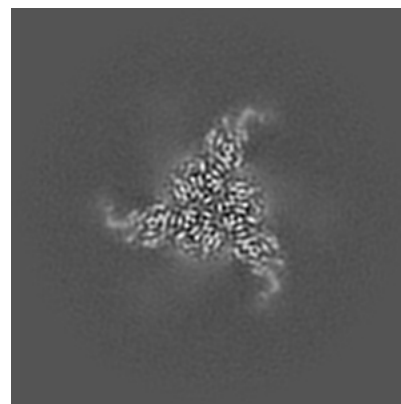
### 6.3.1 Primary map



X Index: 137



Y Index: 134

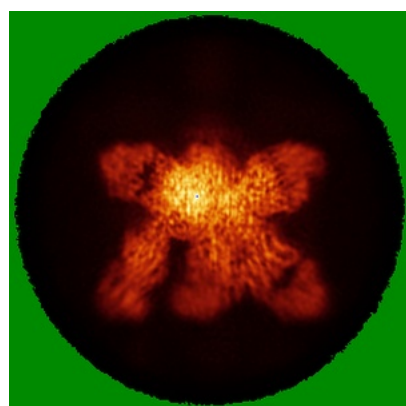


Z Index: 153

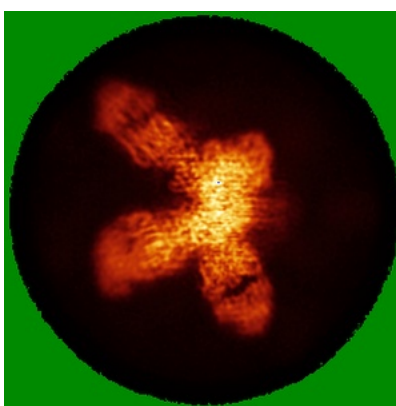
The images above show the largest variance slices of the map in three orthogonal directions.

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

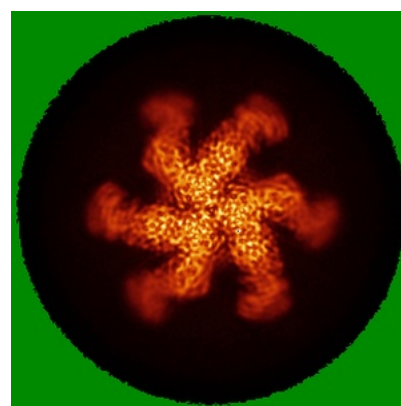
### 6.4.1 Primary map



X



Y

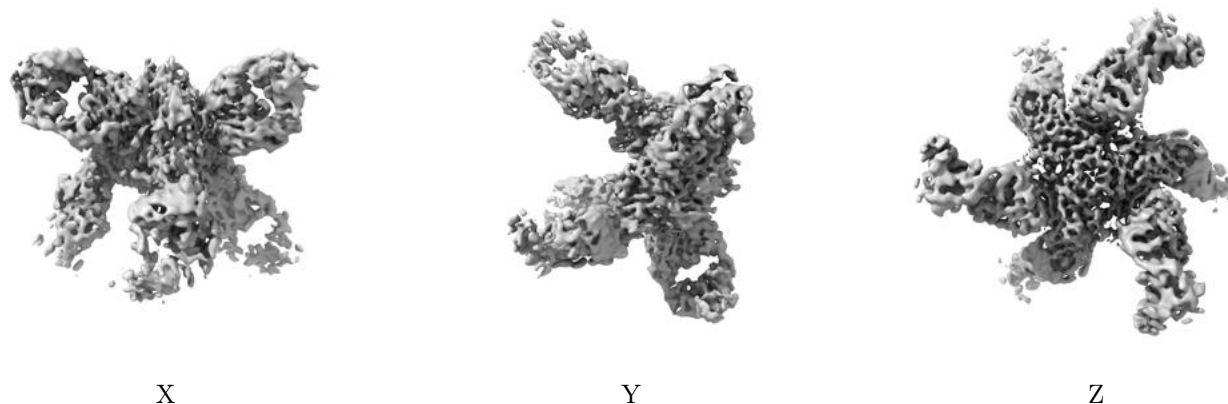


Z

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 0.41. 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](#)

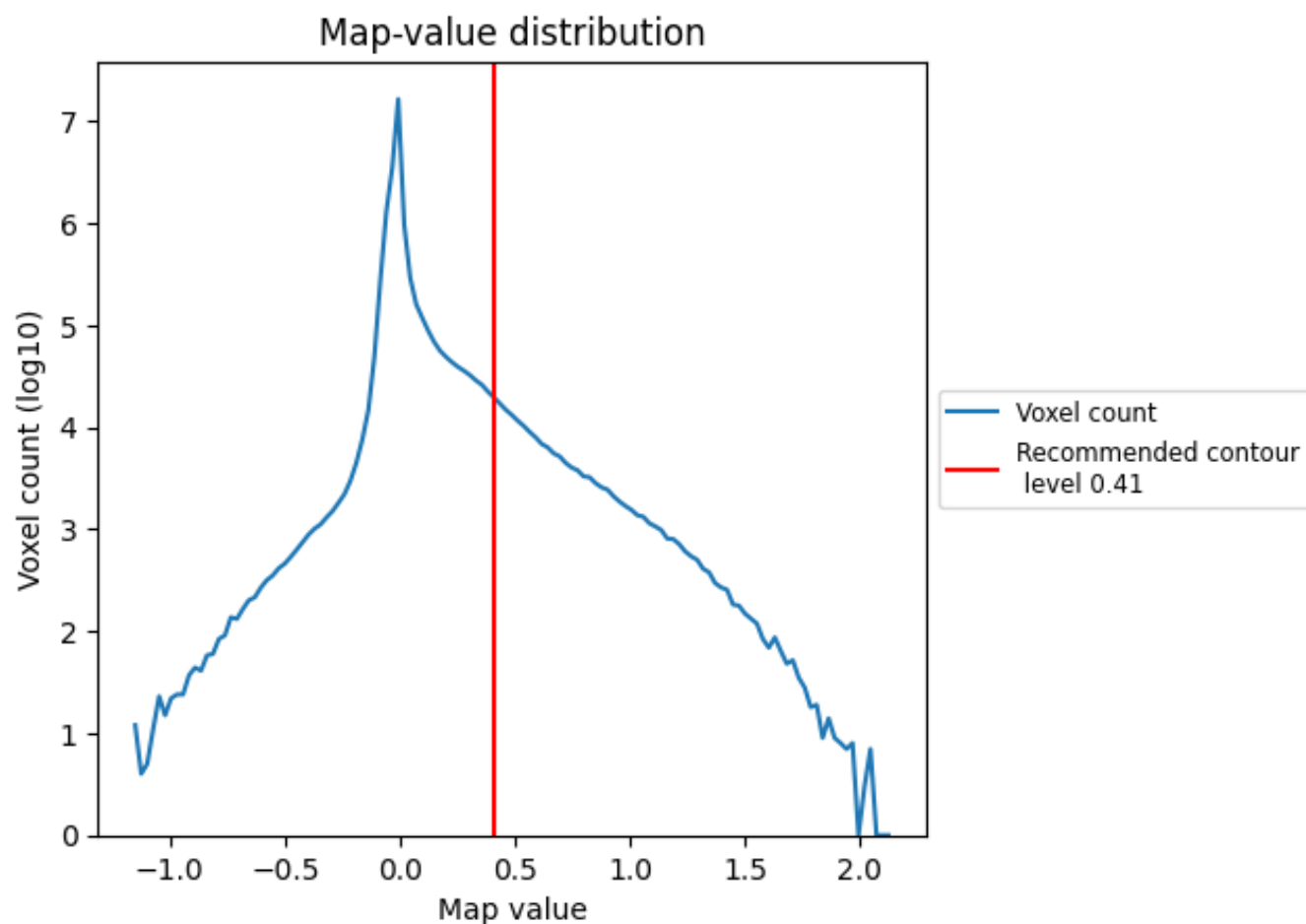
This section was not generated. No masks/segmentation were deposited.



## 7 Map analysis [i](#)

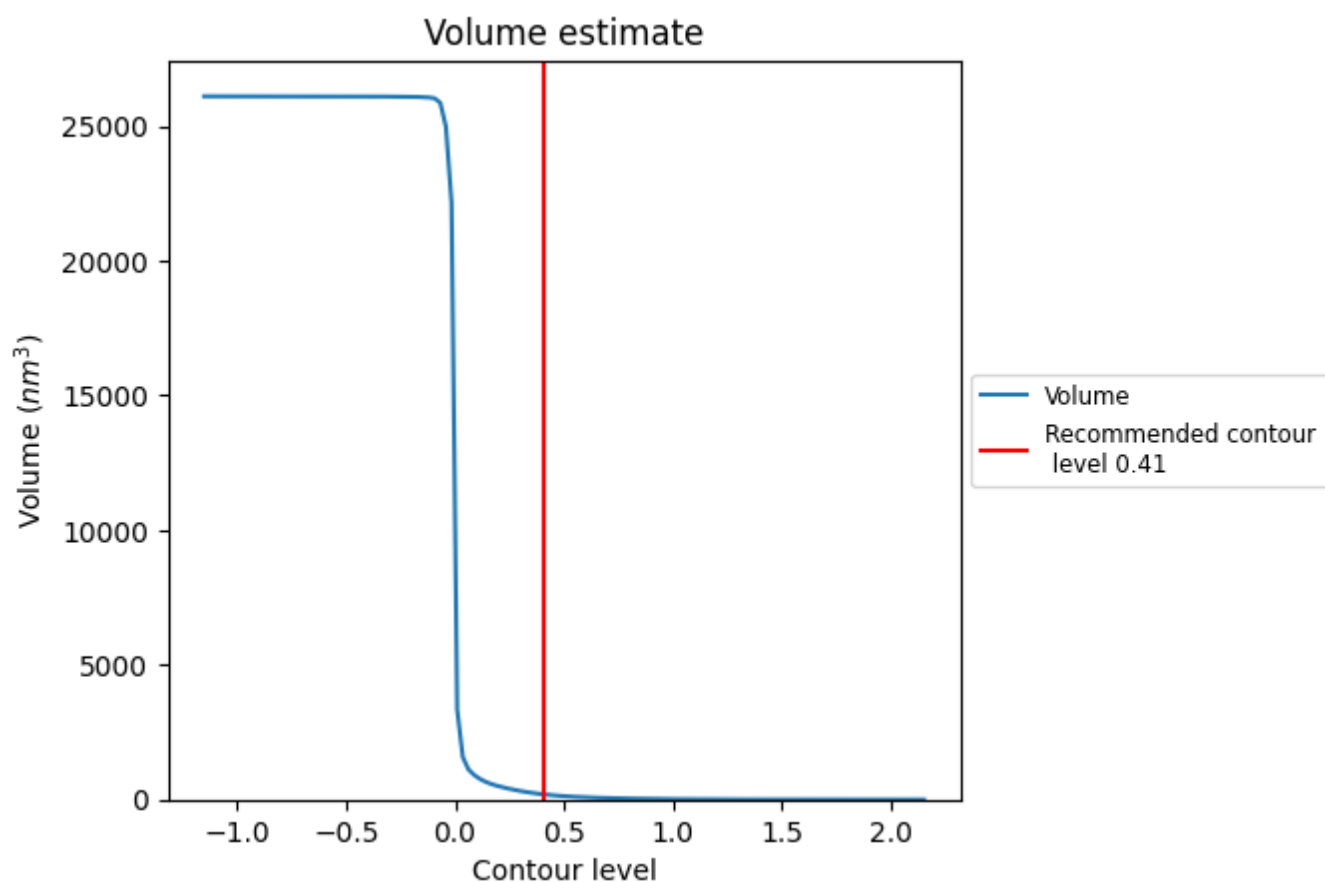
This section contains the results of statistical analysis of the map.

### 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.

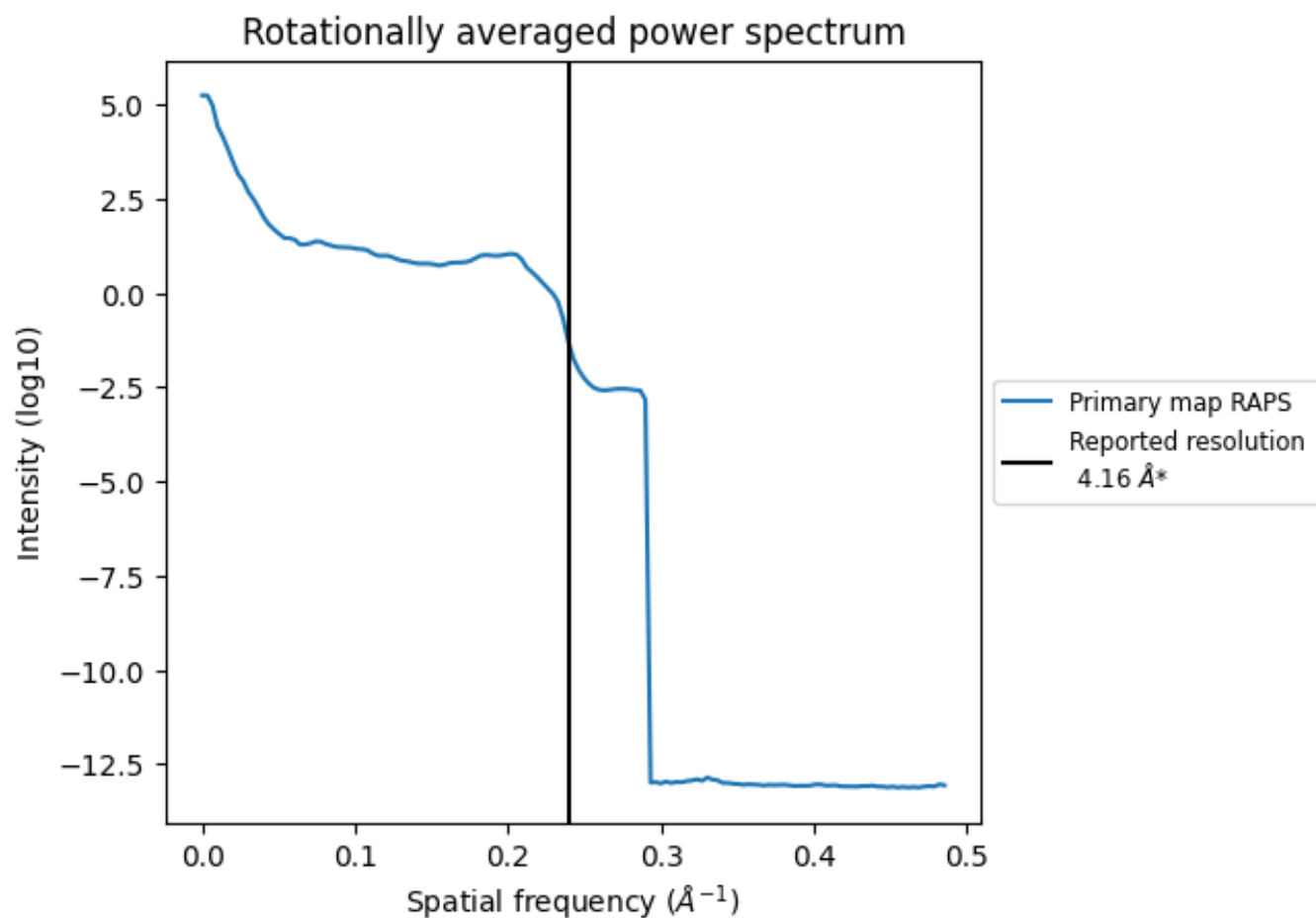
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 190 nm<sup>3</sup>; this corresponds to an approximate mass of 171 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 ⓘ

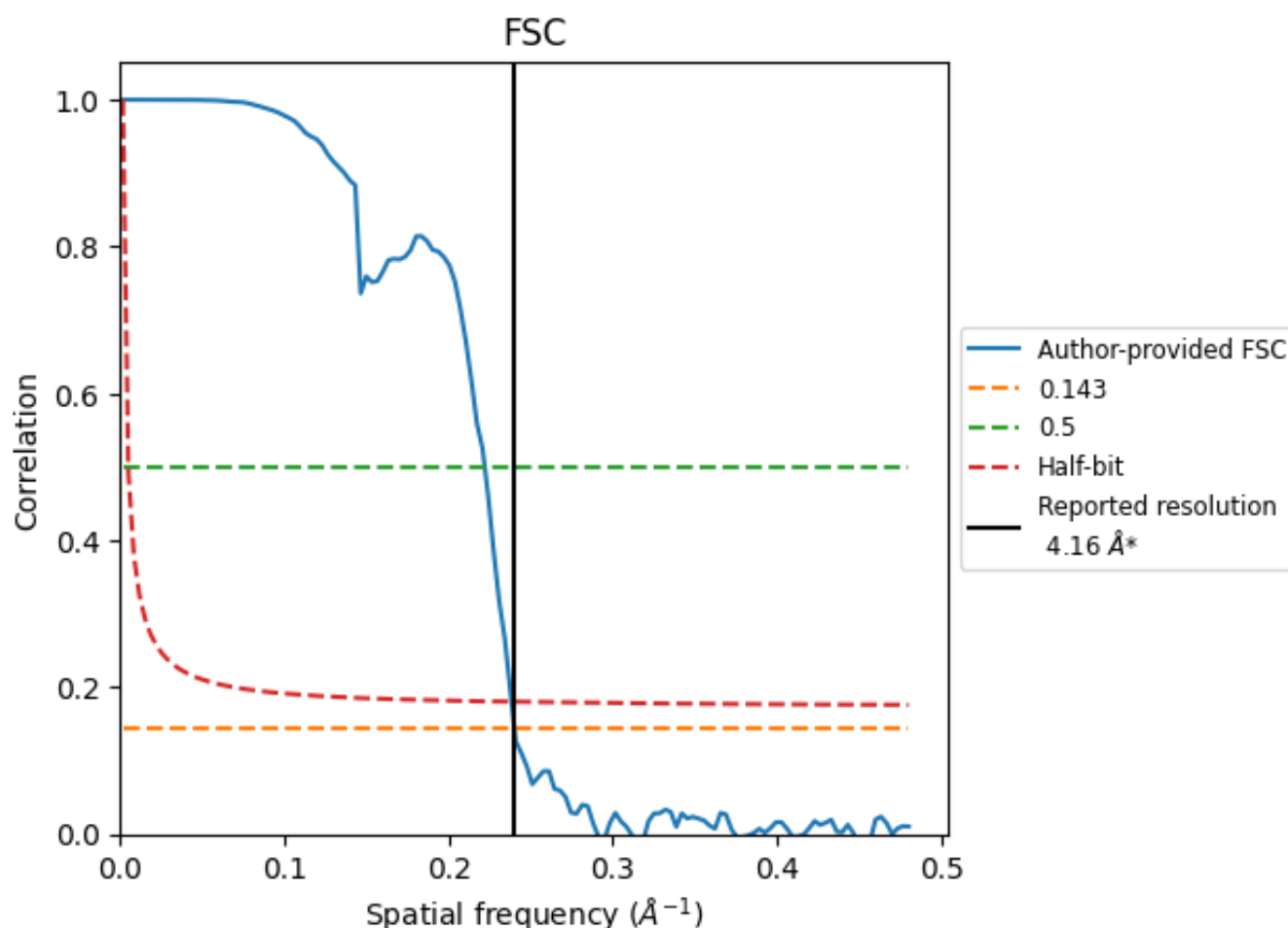


\*Reported resolution corresponds to spatial frequency of 0.240  $\text{\AA}^{-1}$

## 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 0.240 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

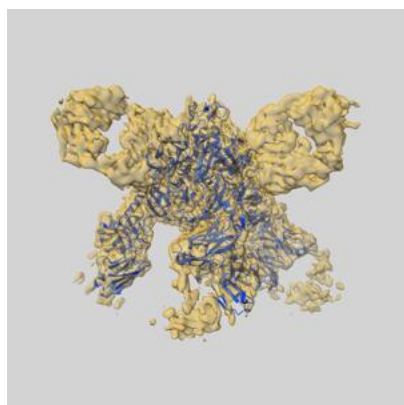
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.16	-	-
Author-provided FSC curve	4.16	4.50	4.19
Unmasked-calculated*	-	-	-

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

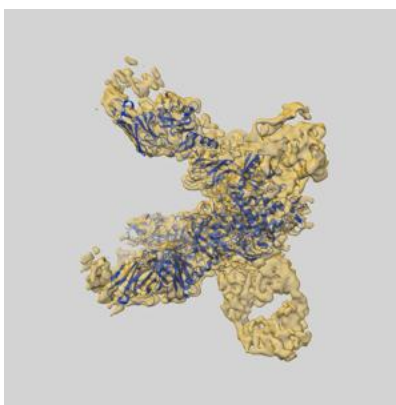
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22841 and PDB model 7KEW. Per-residue inclusion information can be found in section [3](#) on page [12](#).

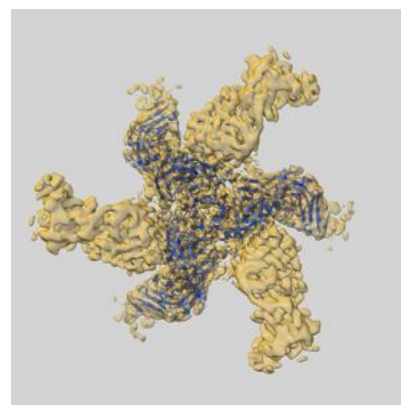
### 9.1 Map-model overlay [i](#)



X



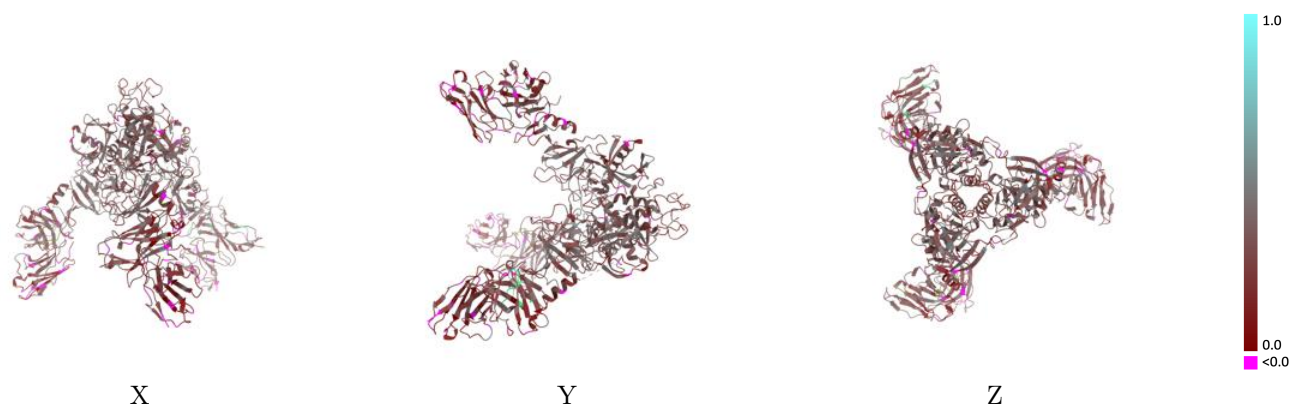
Y



Z

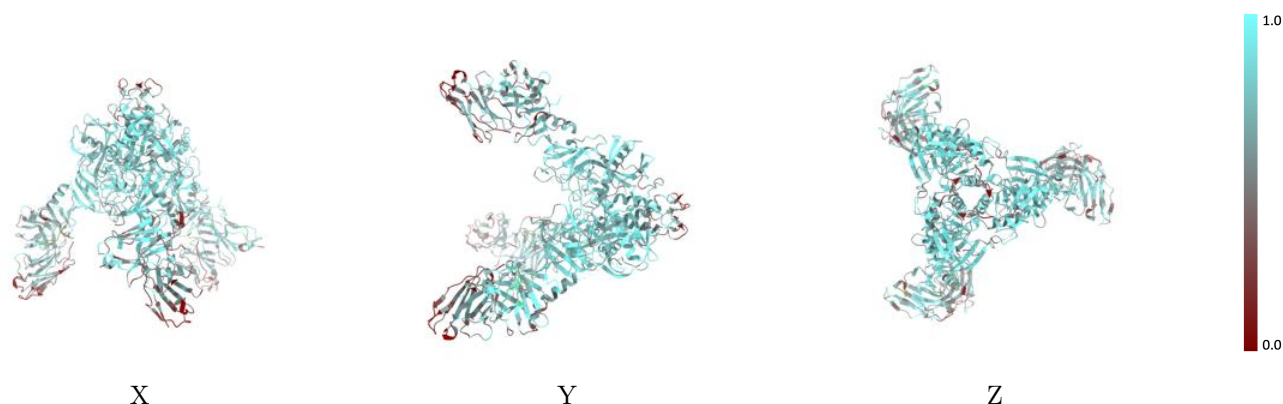
The images above show the 3D surface view of the map at the recommended contour level 0.41 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



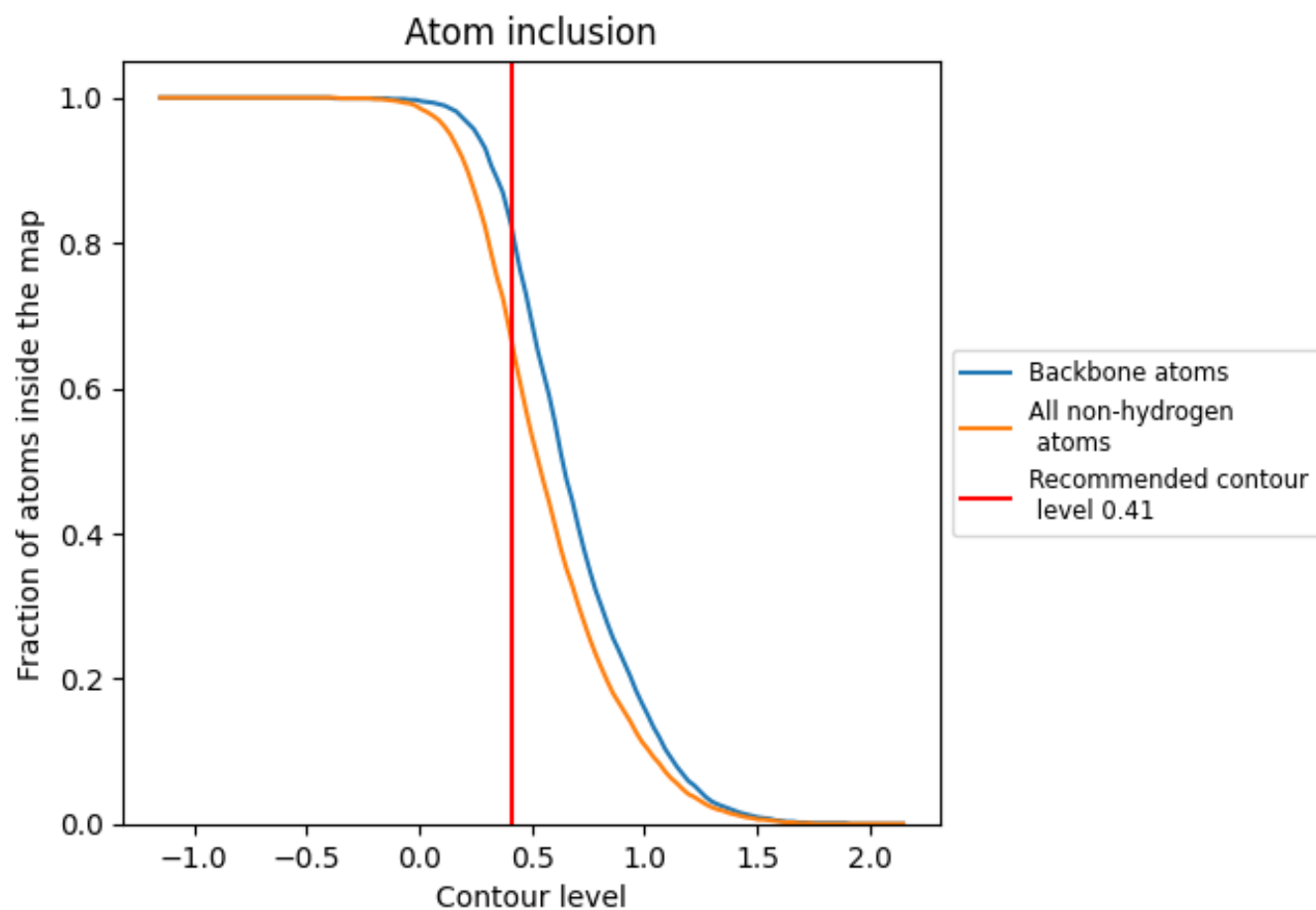
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.41).

## 9.4 Atom inclusion [i](#)









































At the recommended contour level, 82% of all backbone atoms, 67% 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 (0.41) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6660	 0.2680
A	 0.7410	 0.3100
B	 0.7410	 0.3130
C	 0.7420	 0.3130
D	 0.7290	 0.2890
E	 0.7330	 0.2930
F	 0.7330	 0.2920
G	 0.6390	 0.2340
H	 0.6390	 0.2280
I	 0.6390	 0.2290
J	 0.4650	 0.1850
K	 0.4690	 0.1850
L	 0.4650	 0.1870
M	 0.5360	 0.2630
N	 0.5000	 0.2730
O	 0.7210	 0.3630
P	 0.7210	 0.3650
Q	 0.5000	 0.2550
R	 0.7380	 0.3680

