



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 11:23 AM BST

PDB ID : 6K7F  
Title : Crystal structure of MBPholo-Tim21 fusion protein with a 17-residue helical linker  
Authors : Bala, S.; Shimada, A.; Kohda, D.  
Deposited on : 2019-06-07  
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

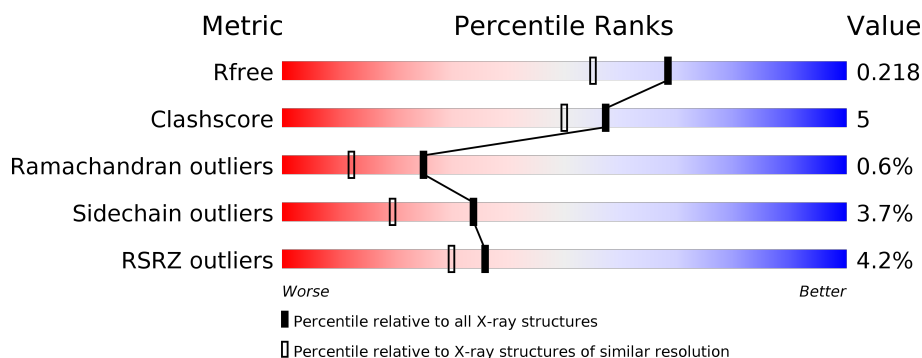
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	499	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>11%</div> <div></div> </div> </div>
2	B	2	<div> <div>100%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4264 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Maltose/maltodextrin-binding periplasmic protein, Mitochondrial import inner membrane translocase subunit TIM21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	499	Total	C	N	O	S	0	0	0
			3922	2511	658	742	11			

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AEX9
A	313	VAL	ALA	engineered mutation	UNP P0AEX9
A	370	GLU	-	linker	UNP P0AEX9
A	371	ALA	-	linker	UNP P0AEX9
A	372	ALA	-	linker	UNP P0AEX9
A	373	ALA	-	linker	UNP P0AEX9
A	374	LYS	-	linker	UNP P0AEX9
A	375	GLU	-	linker	UNP P0AEX9
A	376	ALA	-	linker	UNP P0AEX9
A	377	ALA	-	linker	UNP P0AEX9
A	378	ALA	-	linker	UNP P0AEX9
A	379	LYS	-	linker	UNP P0AEX9
A	380	GLU	-	linker	UNP P0AEX9
A	381	ALA	-	linker	UNP P0AEX9
A	382	ALA	-	linker	UNP P0AEX9
A	383	ALA	-	linker	UNP P0AEX9
A	384	LYS	-	linker	UNP P0AEX9
A	385	ALA	-	linker	UNP P0AEX9
A	386	ALA	-	linker	UNP P0AEX9

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	0	0	0
			23	12	11			

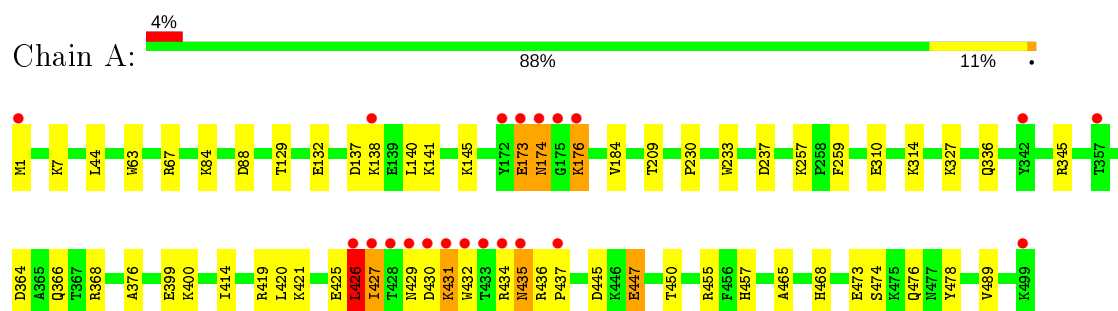
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	319	Total	O	0	0
			319	319		

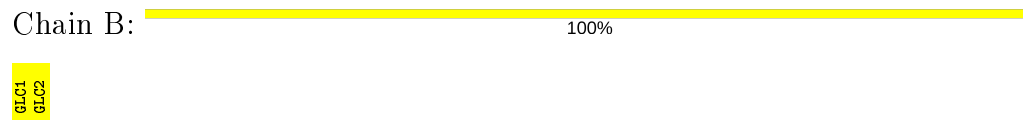
### 3 Residue-property plots

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 electron 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 dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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: Maltose/maltodextrin-binding periplasmic protein,Mitochondrial import inner membrane translocase subunit TIM21



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	40.97Å 69.41Å 171.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.28 – 1.80 35.28 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.28-1.80) 99.9 (35.28-1.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.65 (at 1.81Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, $R_{free}$	0.174 , 0.216 0.176 , 0.218	Depositor DCC
$R_{free}$ test set	2314 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.5	Xtriage
Anisotropy	0.545	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4264	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GLC

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	0.48	0/4009	0.63	2/5423 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	44	LEU	CA-CB-CG	5.86	128.77	115.30
1	A	44	LEU	CB-CG-CD1	-5.01	102.48	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	427	ILE	Peptide

### 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	3922	0	3924	39	0
2	B	23	0	21	0	0
3	A	319	0	0	6	1
All	All	4264	0	3945	39	1

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 5.

All (39) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LYS:NZ	3:A:601:HOH:O	1.76	1.16
1:A:426:LEU:HB3	1:A:427:ILE:HB	1.49	0.91
1:A:209:THR:OG1	3:A:602:HOH:O	1.94	0.86
1:A:476:GLN:N	3:A:604:HOH:O	2.12	0.81
1:A:450:THR:HB	1:A:473:GLU:HG3	1.67	0.77
1:A:426:LEU:HG	1:A:457:HIS:NE2	2.03	0.73
1:A:478:TYR:OH	3:A:603:HOH:O	2.10	0.69
1:A:237:ASP:OD2	3:A:605:HOH:O	2.15	0.62
1:A:173:GLU:OE2	1:A:174:ASN:N	2.32	0.59
1:A:399:GLU:HG2	1:A:420:LEU:HD12	1.86	0.58
1:A:445:ASP:OD1	1:A:447:GLU:HG3	2.05	0.56
1:A:129:THR:OG1	1:A:132:GLU:HG3	2.06	0.56
1:A:336:GLN:HE21	1:A:376:ALA:HB1	1.71	0.54
1:A:426:LEU:HB3	1:A:427:ILE:CB	2.32	0.54
1:A:174:ASN:O	1:A:176:LYS:NZ	2.38	0.53
1:A:431:LYS:HB3	1:A:435:ASN:O	2.11	0.51
1:A:426:LEU:HA	1:A:436:ARG:NH1	2.27	0.50
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.47	0.49
1:A:140:LEU:HB3	1:A:145:LYS:HB2	1.96	0.47
1:A:430:ASP:HA	1:A:432:TRP:CE2	2.50	0.47
1:A:431:LYS:HG3	1:A:437:PRO:HD3	1.97	0.45
1:A:425:GLU:HA	1:A:426:LEU:HB2	1.99	0.45
1:A:364:ASP:O	1:A:368:ARG:HG3	2.18	0.44
1:A:84:LYS:HE2	1:A:88:ASP:OD1	2.18	0.44
1:A:419:ARG:NH1	3:A:606:HOH:O	2.24	0.43
1:A:336:GLN:HE21	1:A:376:ALA:CB	2.32	0.43
1:A:137:ASP:OD2	1:A:141:LYS:HE3	2.18	0.43
1:A:336:GLN:NE2	1:A:376:ALA:HB1	2.32	0.42
1:A:421:LYS:HB3	1:A:421:LYS:HE3	1.82	0.42
1:A:426:LEU:H	1:A:427:ILE:C	2.22	0.42
1:A:425:GLU:CA	1:A:426:LEU:HB2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:O	1:A:141:LYS:HG2	2.20	0.41
1:A:465:ALA:HB2	1:A:489:VAL:HG22	2.01	0.41
1:A:257:LYS:HG2	1:A:327:LYS:O	2.19	0.41
1:A:7:LYS:HE3	1:A:7:LYS:HB3	1.76	0.41
1:A:184:VAL:HG22	1:A:366:GLN:OE1	2.21	0.41
1:A:455:ARG:HD3	1:A:468:HIS:CD2	2.56	0.41
1:A:430:ASP:HA	1:A:432:TRP:NE1	2.36	0.40
1:A:63:TRP:CD1	1:A:67:ARG:HG3	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:603:HOH:O	3:A:707:HOH:O[3_654]	2.11	0.09

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	497/499 (100%)	481 (97%)	13 (3%)	3 (1%)	25	12

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	429	ASN
1	A	434	ARG
1	A	426	LEU

### 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 X-ray entries followed by that with respect to entries of similar resolution.

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	407/407 (100%)	392 (96%)	15 (4%)	34	19

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	173	GLU
1	A	174	ASN
1	A	176	LYS
1	A	259	PHE
1	A	310	GLU
1	A	314	LYS
1	A	345	ARG
1	A	400	LYS
1	A	414	ILE
1	A	426	LEU
1	A	431	LYS
1	A	435	ASN
1	A	447	GLU
1	A	474	SER

Some 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 ⓘ

2 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
2	GLC	B	1	2	12,12,12	0.68	0	17,17,17	1.06	3 (17%)
2	GLC	B	2	2	11,11,12	0.85	0	15,15,17	1.05	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
2	GLC	B	1	2	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	GLC	C1-O5-C5	2.75	115.92	112.19
2	B	1	GLC	O2-C2-C3	-2.11	105.47	110.35
2	B	1	GLC	C6-C5-C4	-2.08	108.13	113.00
2	B	1	GLC	O4-C4-C3	-2.02	105.68	110.35

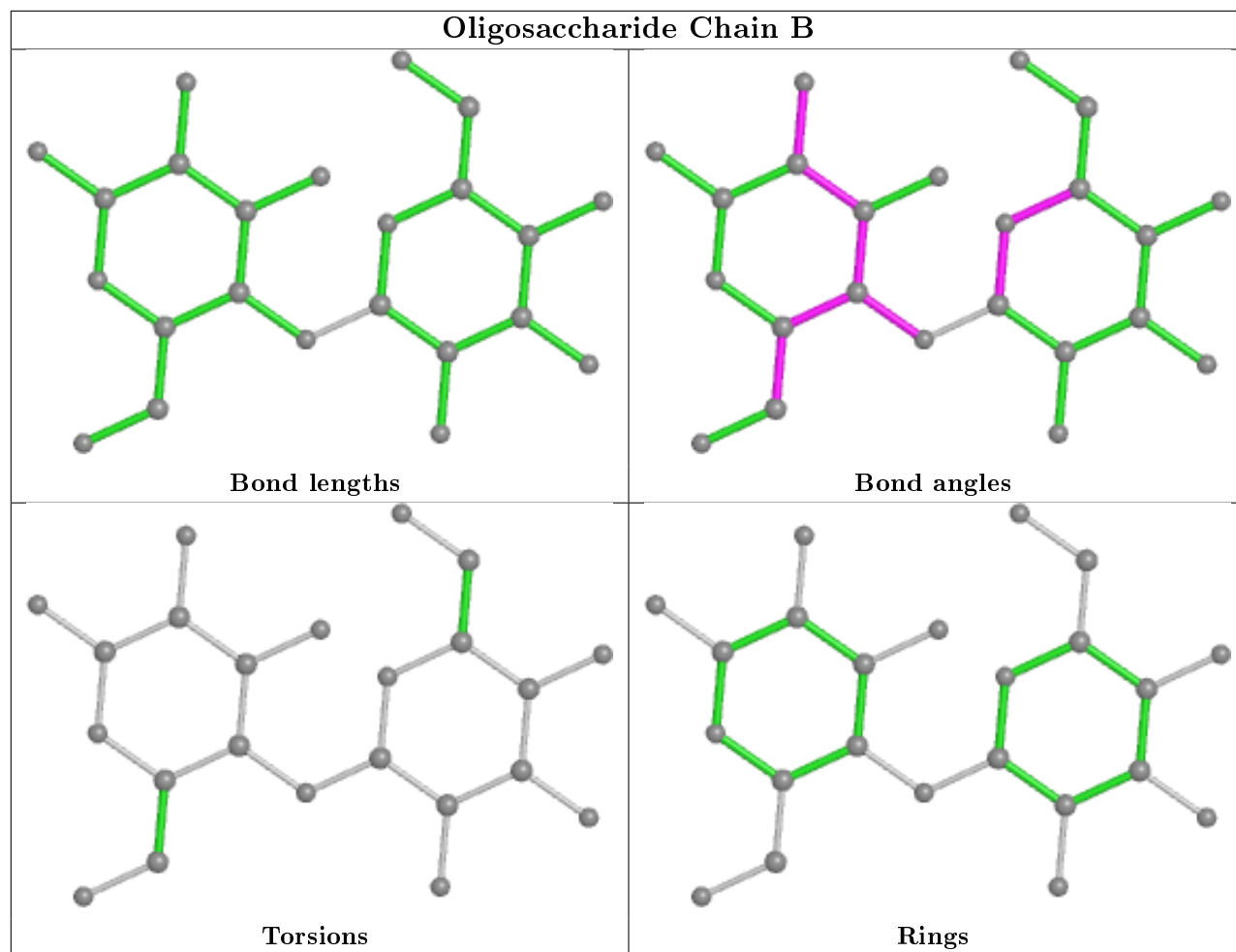
There are no chirality outliers.

There are no torsion outliers.

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](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	499/499 (100%)	0.01	21 (4%) 36 30	18, 36, 66, 111	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	433	THR	8.2
1	A	430	ASP	7.9
1	A	429	ASN	7.5
1	A	427	ILE	7.4
1	A	173	GLU	5.7
1	A	432	TRP	5.2
1	A	426	LEU	5.0
1	A	174	ASN	4.3
1	A	428	THR	3.8
1	A	172	TYR	3.6
1	A	176	LYS	3.6
1	A	434	ARG	3.2
1	A	435	ASN	3.2
1	A	175	GLY	3.1
1	A	431	LYS	3.1
1	A	499	LYS	3.1
1	A	1	MET	2.8
1	A	437	PRO	2.1
1	A	342	TYR	2.1
1	A	357	THR	2.0
1	A	138	LYS	2.0

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

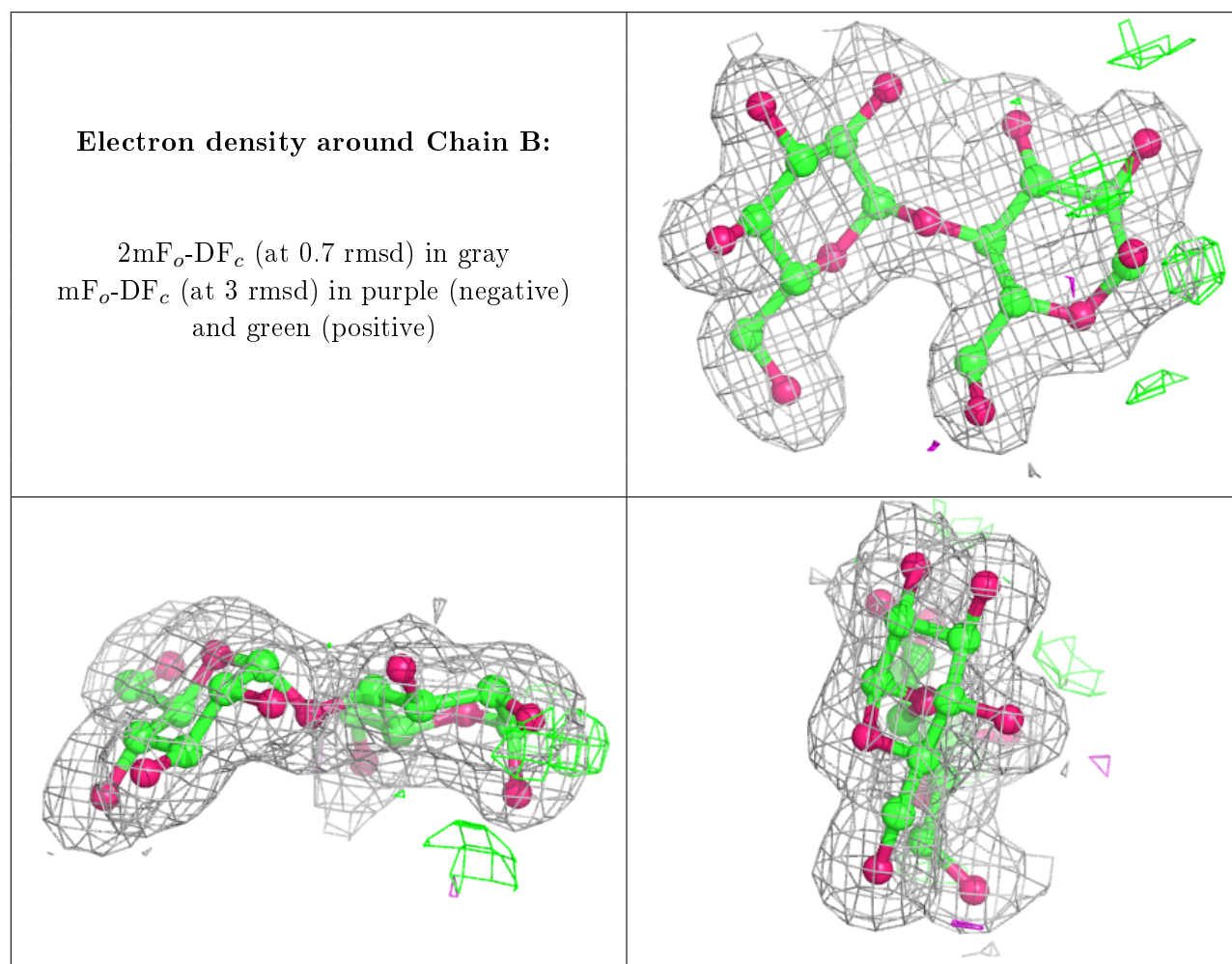
There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	GLC	B	1	12/12	0.95	0.10	18,26,34,34	0
2	GLC	B	2	11/12	0.97	0.11	16,21,23,23	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



### 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.