



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 02:40 AM UTC

PDB ID : 4WEI / pdb\_00004wei  
Title : Crystal structure of the F4 fimbrial adhesin FaeG in complex with lactose  
Authors : Moonens, K.; Van den Broeck, I.; De Kerpel, M.; Deboeck, F.; Raymaekers, H.; Remaut, H.; De Greve, H.  
Deposited on : 2014-09-10  
Resolution : 2.30 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.

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:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

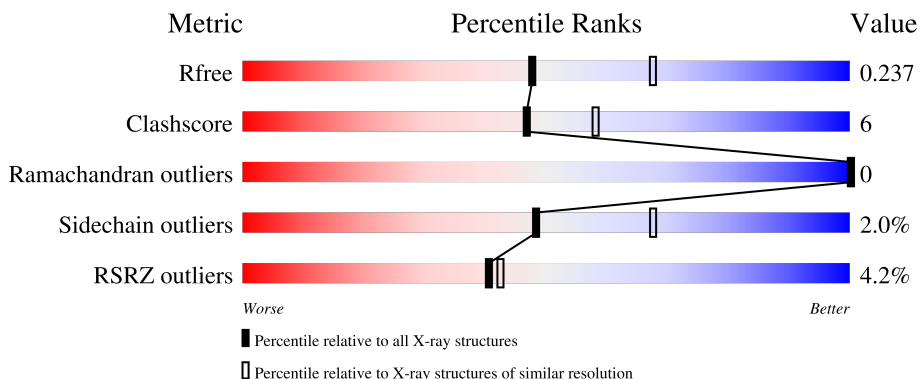
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



The reported resolution of this entry is 2.30 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))
$R_{free}$	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 of 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	277	 4% 83% 9% • 6%
2	B	2	 50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	B	1	X	-	-	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2052 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 K88 fimbrial protein AD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	259	1921	1206	324	387	4	0	3	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	TRP	-	expression tag	UNP P14191
A	10	MET	-	expression tag	UNP P14191
A	11	THR	-	expression tag	UNP P14191
A	12	GLY	-	expression tag	UNP P14191
A	13	HIS	-	expression tag	UNP P14191
A	14	HIS	-	expression tag	UNP P14191
A	15	HIS	-	expression tag	UNP P14191
A	16	HIS	-	expression tag	UNP P14191
A	17	HIS	-	expression tag	UNP P14191
A	18	HIS	-	expression tag	UNP P14191
A	38	SER	ASN	conflict	UNP P14191
A	265	ASP	-	expression tag	UNP P14191
A	266	ASN	-	expression tag	UNP P14191
A	267	LYS	-	expression tag	UNP P14191
A	268	GLN	-	expression tag	UNP P14191
A	269	MET	-	expression tag	UNP P14191
A	270	THR	-	expression tag	UNP P14191
A	271	GLY	-	expression tag	UNP P14191
A	272	ASP	-	expression tag	UNP P14191
A	273	PHE	-	expression tag	UNP P14191
A	274	ASN	-	expression tag	UNP P14191
A	275	GLY	-	expression tag	UNP P14191
A	276	SER	-	expression tag	UNP P14191
A	277	VAL	-	expression tag	UNP P14191
A	278	ASP	-	expression tag	UNP P14191
A	279	ILE	-	expression tag	UNP P14191
A	280	GLY	-	expression tag	UNP P14191

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Chain	Residue	Modelled	Actual	Comment	Reference
A	281	GLY	-	expression tag	UNP P14191
A	282	SER	-	expression tag	UNP P14191
A	283	ILE	-	expression tag	UNP P14191
A	284	THR	-	expression tag	UNP P14191
A	285	ALA	-	expression tag	UNP P14191

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-alpha-D-glucopyranoside.



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

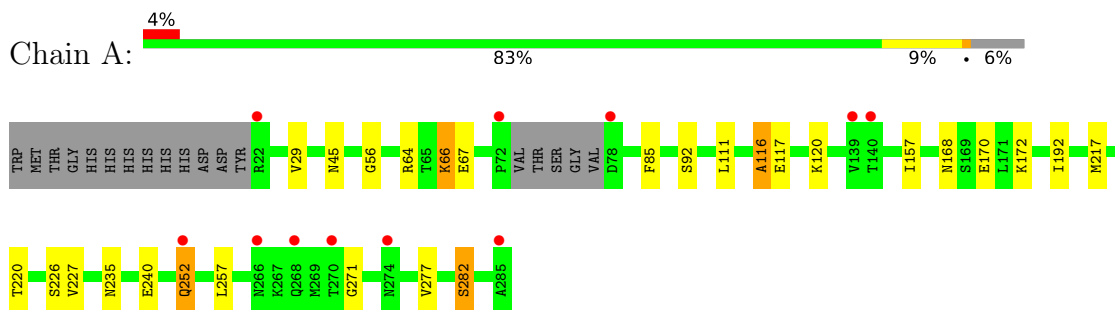
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	108	108	108	0	0

### 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 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: K88 fimbrial protein AD



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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.19Å 93.19Å 110.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.67 – 2.30 47.67 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.67-2.30) 99.9 (47.67-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 2.29Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.192 , 0.230 0.206 , 0.237	Depositor DCC
$R_{free}$ test set	1143 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.4	Xtrriage
Anisotropy	0.044	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 28.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2052	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.46% 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 [i](#)

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

Bond lengths and bond angles in the following residue types are not validated in this section: GAL, 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	1.21	2/1953 (0.1%)	1.14	3/2653 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	SER	CA-C	-5.50	1.45	1.52
1	A	282	SER	N-CA	-5.44	1.39	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	LYS	N-CA-C	-5.55	106.34	113.23
1	A	116	ALA	N-CA-C	-5.54	102.08	110.28
1	A	45	ASN	CB-CA-C	-5.35	104.30	112.11

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1921	0	1882	22	0
2	B	23	0	20	1	0
3	A	108	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	2052	0	1902	22	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 6.

All (22) 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:252:GLN:HE21	1:A:282:SER:HB3	1.37	0.90
1:A:252:GLN:NE2	1:A:282:SER:HB3	1.86	0.89
1:A:116:ALA:O	1:A:117:GLU:HB2	1.79	0.80
1:A:252:GLN:HE21	1:A:282:SER:CB	1.96	0.78
1:A:252:GLN:NE2	1:A:282:SER:CB	2.48	0.75
1:A:271:GLY:O	3:A:487:HOH:O	2.03	0.74
1:A:116:ALA:O	1:A:117:GLU:CB	2.36	0.73
1:A:64:ARG:NH1	3:A:401:HOH:O	2.19	0.72
1:A:240:GLU:OE1	3:A:485:HOH:O	2.08	0.71
1:A:56:GLY:H	1:A:235:ASN:HD22	1.44	0.64
1:A:157:ILE:HD12	1:A:217:MET:HE3	1.84	0.58
1:A:157:ILE:CD1	1:A:217:MET:HE3	2.41	0.50
1:A:252:GLN:NE2	1:A:282:SER:HB2	2.27	0.50
1:A:29:VAL:HB	1:A:277:VAL:HG22	1.94	0.50
1:A:85:PHE:CD1	1:A:257[B]:LEU:HD11	2.48	0.48
1:A:168:ASN:HD22	1:A:172:LYS:NZ	2.12	0.48
1:A:192:ILE:HG23	1:A:226:SER:HB3	1.96	0.47
1:A:56:GLY:H	1:A:235:ASN:ND2	2.12	0.47
1:A:56:GLY:N	1:A:235:ASN:HD22	2.14	0.46
1:A:66:LYS:C	1:A:67:GLU:HG3	2.45	0.41
1:A:111:LEU:HD13	1:A:257[A]:LEU:HD13	2.02	0.41
1:A:170:GLU:OE2	2:B:2:GAL:O3	2.37	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

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

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	258/277 (93%)	255 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	201/215 (94%)	197 (98%)	4 (2%)	48	67

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	LYS
1	A	220	THR
1	A	227	VAL
1	A	252	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	ASN
1	A	200	ASN
1	A	235	ASN
1	A	252	GLN
1	A	258	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	1.25	1 (8%)	17,17,17	3.63	7 (41%)
2	GAL	B	2	2	11,11,12	0.84	0	15,15,17	1.42	2 (13%)

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	1/1/5/5	1/2/22/22	0/1/1/1
2	GAL	B	2	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	GLC	O4-C4	-3.47	1.34	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	GLC	O4-C4-C5	12.99	141.32	109.32
2	B	1	GLC	C1-C2-C3	-4.51	101.16	110.36
2	B	2	GAL	C2-C3-C4	3.46	116.95	110.86
2	B	2	GAL	O5-C5-C6	2.71	112.94	107.66
2	B	1	GLC	O3-C3-C4	2.59	116.47	110.38
2	B	1	GLC	O2-C2-C1	2.56	115.15	109.25
2	B	1	GLC	C4-C3-C2	-2.46	106.52	110.83
2	B	1	GLC	C1-O5-C5	2.37	118.24	113.65
2	B	1	GLC	O5-C5-C6	2.31	112.17	106.44

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1	GLC	C4

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	GLC	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	GAL	1	0

## 5.5 Carbohydrates [i](#)

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	1.25	1 (8%)	17,17,17	3.63	7 (41%)
2	GAL	B	2	2	11,11,12	0.84	0	15,15,17	1.42	2 (13%)

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	1/1/5/5	1/2/22/22	0/1/1/1
2	GAL	B	2	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	GLC	O4-C4	-3.47	1.34	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	GLC	O4-C4-C5	12.99	141.32	109.32
2	B	1	GLC	C1-C2-C3	-4.51	101.16	110.36
2	B	2	GAL	C2-C3-C4	3.46	116.95	110.86
2	B	2	GAL	O5-C5-C6	2.71	112.94	107.66
2	B	1	GLC	O3-C3-C4	2.59	116.47	110.38
2	B	1	GLC	O2-C2-C1	2.56	115.15	109.25
2	B	1	GLC	C4-C3-C2	-2.46	106.52	110.83
2	B	1	GLC	C1-O5-C5	2.37	118.24	113.65
2	B	1	GLC	O5-C5-C6	2.31	112.17	106.44

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	1	GLC	C4

All (1) torsion outliers are listed below:

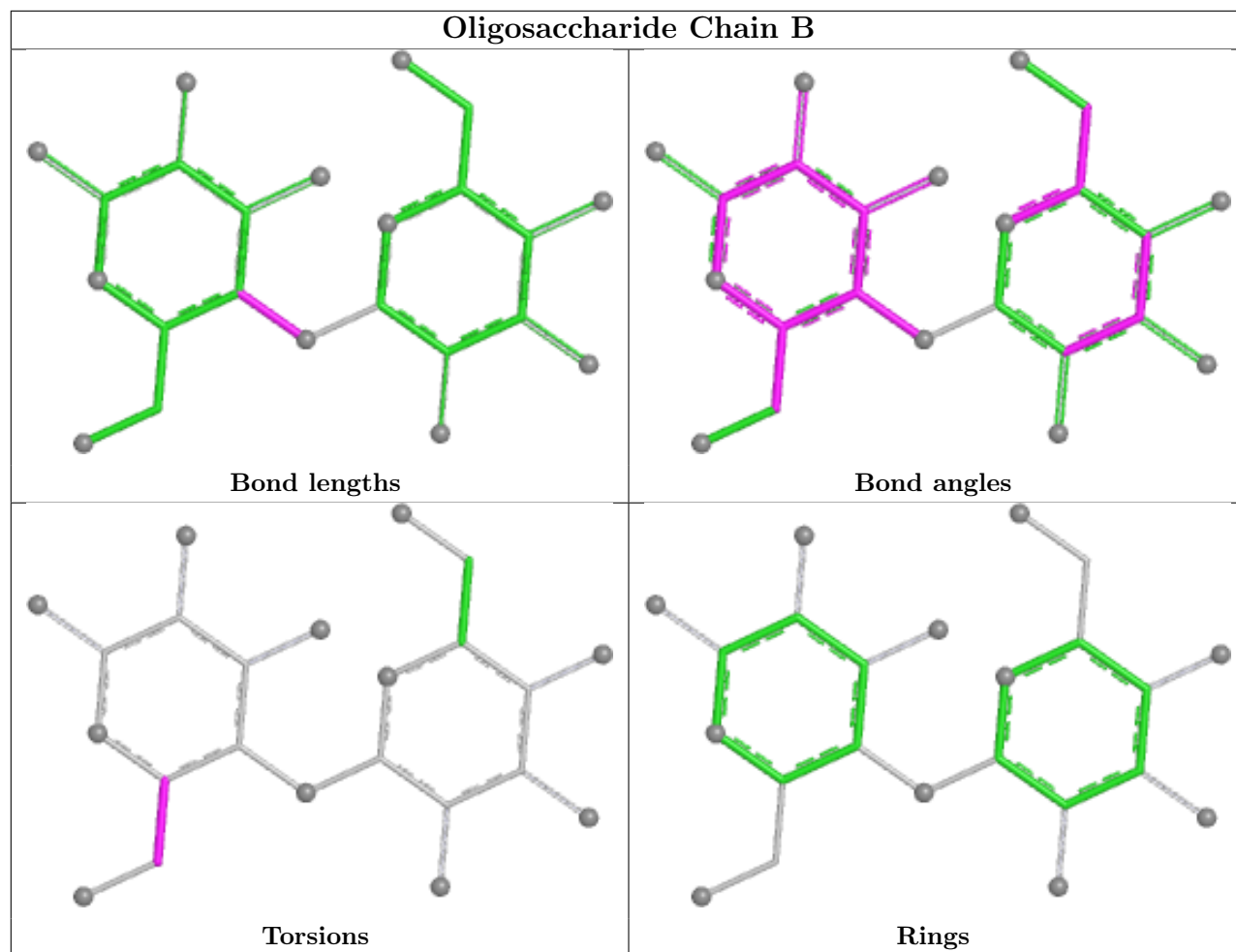
Mol	Chain	Res	Type	Atoms
2	B	1	GLC	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	GAL	1	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](#)

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	259/277 (93%)	0.04	11 (4%) 40 42	18, 41, 62, 83	3 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	22	ARG	4.6
1	A	72	PRO	4.0
1	A	252	GLN	4.0
1	A	268	GLN	3.7
1	A	270	THR	3.1
1	A	140	THR	3.0
1	A	285	ALA	3.0
1	A	78	ASP	2.8
1	A	274	ASN	2.2
1	A	139	VAL	2.1
1	A	266	ASN	2.1

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

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.

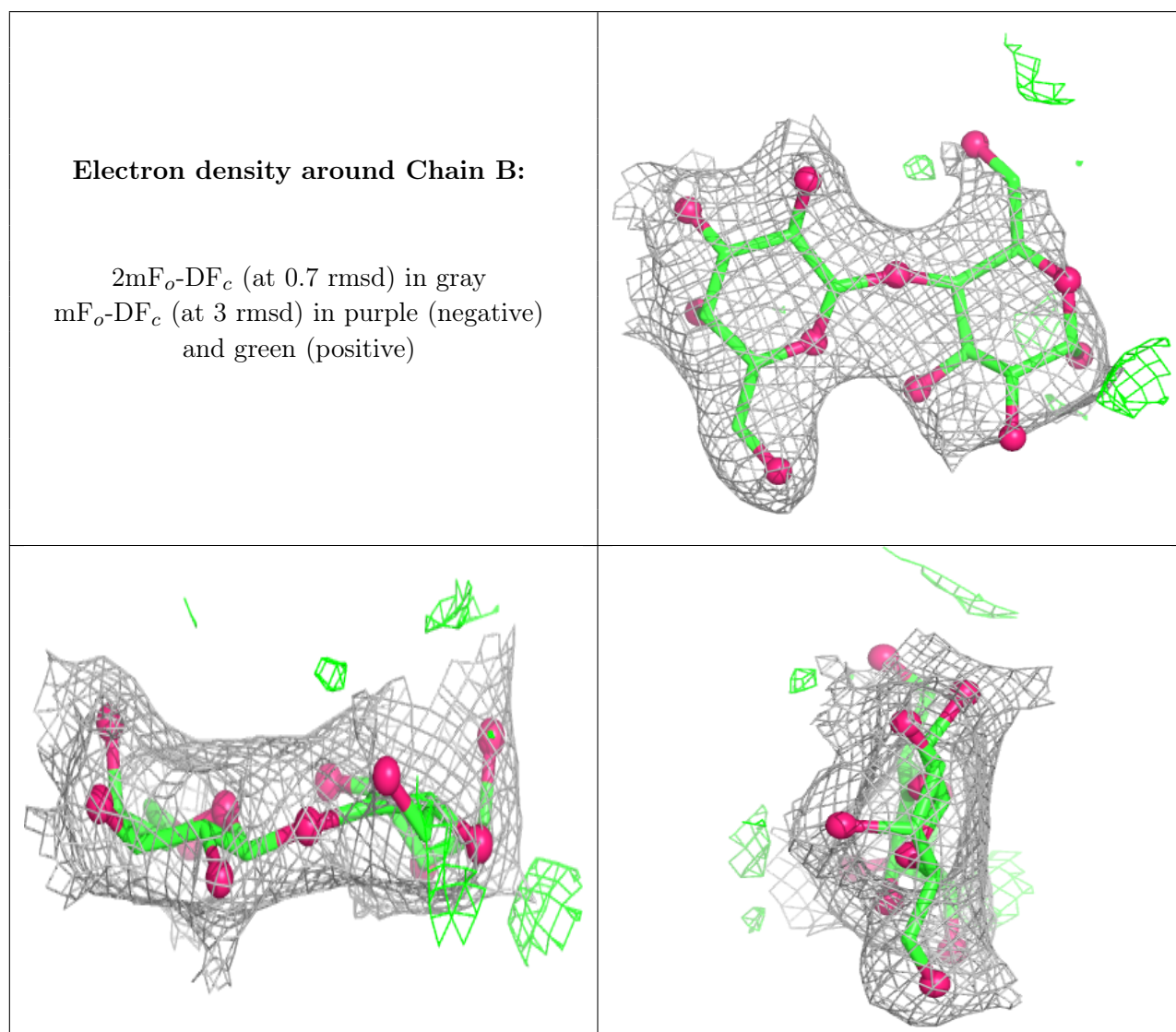
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	GLC	B	1	12/12	0.83	0.20	67,87,99,101	0
2	GAL	B	2	11/12	0.96	0.07	38,49,55,56	0

### 6.3 Carbohydrates [i](#)

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	0.83	0.20	67,87,99,101	0
2	GAL	B	2	11/12	0.96	0.07	38,49,55,56	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

There are no such residues in this entry.