



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 04:49 PM EDT

PDB ID : 3VLW  
Title : Crystal structure of Sphingomonas sp. A1 alginate-binding protein AlgQ1 in complex with mannuronate-guluronate disaccharide  
Authors : Nishitani, Y.; Maruyama, Y.; Itoh, T.; Mikami, B.; Hashimoto, W.; Murata, K.  
Deposited on : 2011-12-05  
Resolution : 2.00 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
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)	:	2.36.2

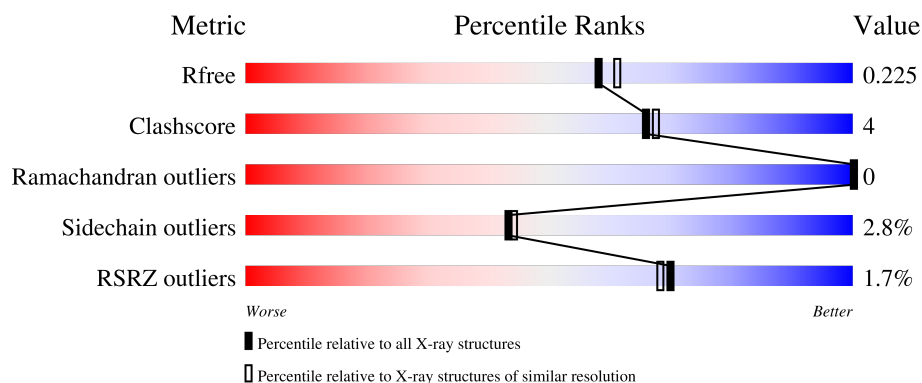
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

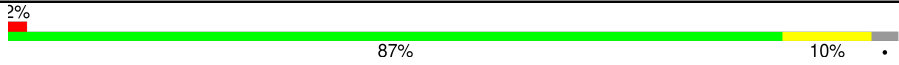



The reported resolution of this entry is 2.00 Å.

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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	502	
1	B	502	
2	C	2	
2	D	2	

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
4	GOL	B	516	-	-	-	X

## 2 Entry composition [i](#)

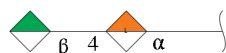
There are 5 unique types of molecules in this entry. The entry contains 8758 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 AlgQ1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	5	0
			4045	2612	688	731	14			
1	B	489	Total	C	N	O	S	6	13	0
			4095	2650	696	735	14			

- Molecule 2 is an oligosaccharide called beta-D-mannopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			25	12	13			
2	D	2	Total	C	O	0	0	0
			25	12	13			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	235	Total	O	0	0
			235	235		
5	B	223	Total	O	0	0
			223	223		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.76Å 67.53Å 90.69Å 90.00° 93.33° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 45.27 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (50.00-2.00) 99.6 (45.27-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.55 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.190 , 0.223 0.192 , 0.225	Depositor DCC
$R_{free}$ test set	3291 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.5	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 51.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8758	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 65.90 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.5894e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: CA, GOL, BEM, LGU

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	0.35	0/4178	0.48	0/5668
1	B	0.37	0/4249	0.49	0/5760
All	All	0.36	0/8427	0.49	0/11428

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [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	4045	0	3960	29	0
1	B	4095	0	4045	36	0
2	C	25	0	15	0	0
2	D	25	0	15	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	42	0	56	1	0
4	B	66	0	88	2	0
5	A	235	0	0	0	0
5	B	223	0	0	2	0
All	All	8758	0	8179	66	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (66) 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:B:253[B]:ARG:HH11	1:B:253[B]:ARG:CG	1.68	1.06
1:B:253[B]:ARG:HH11	1:B:253[B]:ARG:HG3	1.27	0.98
1:B:253[B]:ARG:HH11	1:B:253[B]:ARG:CB	1.81	0.93
1:B:253[B]:ARG:NH1	1:B:253[B]:ARG:HB3	1.88	0.89
1:A:93:ASN:HD21	1:A:124:ILE:H	1.21	0.83
1:B:93:ASN:HD21	1:B:124:ILE:H	1.26	0.83
1:A:110:HIS:O	1:A:113[A]:VAL:HG12	1.82	0.80
1:B:155:GLN:HE21	1:B:303:ARG:HH22	1.31	0.77
1:B:253[B]:ARG:HG3	1:B:253[B]:ARG:NH1	1.97	0.76
1:B:7:THR:HG22	1:B:328:GLU:HG2	1.70	0.73
1:B:185:ASN:HD22	1:B:187:ASP:H	1.36	0.73
1:A:185:ASN:HD22	1:A:187:ASP:H	1.36	0.72
1:B:93:ASN:H	1:B:93:ASN:HD22	1.37	0.72
1:A:106:PHE:CE2	1:A:113[A]:VAL:HG11	2.29	0.68
1:A:93:ASN:HD22	1:A:93:ASN:H	1.42	0.67
1:A:155:GLN:HE21	1:A:303:ARG:HH22	1.43	0.66
1:A:93:ASN:ND2	1:A:124:ILE:H	1.96	0.61
1:B:104[A]:LYS:NZ	1:B:104[A]:LYS:HB3	2.15	0.61
1:B:253[B]:ARG:CB	1:B:253[B]:ARG:NH1	2.49	0.60
1:A:196:ASN:HD21	1:A:306:GLU:HB2	1.68	0.57
1:B:349:VAL:H	1:B:353:THR:HG22	1.69	0.57
1:B:93:ASN:ND2	1:B:124:ILE:H	2.01	0.54
1:A:213:VAL:HG22	1:A:218:ILE:HD13	1.89	0.54
1:B:7:THR:HG22	1:B:328:GLU:CG	2.38	0.54
1:B:196:ASN:HD21	1:B:306:GLU:HB2	1.72	0.53
1:A:78:ASP:HA	1:A:431[A]:ARG:HH22	1.74	0.52
1:A:424:VAL:HA	1:A:489:GLN:HE22	1.74	0.51
1:A:2:GLU:HB2	1:A:5:TRP:CD1	2.45	0.51
1:B:7:THR:CG2	1:B:328:GLU:HG2	2.39	0.50
1:B:192:PHE:O	1:B:195:VAL:HG22	2.12	0.50
1:A:204:GLY:HA3	1:A:211:PHE:O	2.12	0.49
1:A:348:GLY:HA3	1:A:353:THR:HG21	1.95	0.48
1:B:104[A]:LYS:HB3	1:B:104[A]:LYS:HZ3	1.78	0.48
1:B:228:LYS:HD3	1:B:463:TRP:CG	2.49	0.48
1:A:347:PHE:O	1:A:353:THR:CG2	2.62	0.47
1:B:348:GLY:HA3	1:B:353:THR:HG21	1.97	0.47
1:B:93:ASN:H	1:B:93:ASN:ND2	2.10	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213[A]:VAL:HG22	1:B:218:ILE:HD13	1.97	0.47
1:B:213[B]:VAL:HG11	4:B:510:GOL:H11	1.96	0.47
1:A:93:ASN:H	1:A:93:ASN:ND2	2.12	0.46
1:A:192:PHE:O	1:A:195:VAL:HG22	2.17	0.45
1:A:349:VAL:H	1:A:353:THR:HG22	1.82	0.45
4:B:507:GOL:H11	5:B:594:HOH:O	2.17	0.45
1:A:79:LYS:HG2	4:A:511:GOL:H32	1.99	0.44
1:A:93:ASN:HD21	1:A:124:ILE:N	2.01	0.44
1:B:353:THR:OG1	1:B:367:VAL:HG21	2.18	0.44
1:B:93:ASN:HD21	1:B:124:ILE:N	2.06	0.43
1:B:104[A]:LYS:NZ	1:B:104[A]:LYS:CB	2.81	0.43
1:A:269:ASP:HB3	1:A:274:THR:OG1	2.19	0.43
1:A:458:ASP:HA	1:A:459:PRO:HD2	1.91	0.43
1:A:199:GLY:HA3	1:A:300:LYS:HG3	2.02	0.42
1:B:48:ASN:ND2	1:B:50:ALA:H	2.18	0.42
1:B:119:ALA:HB1	1:B:120:PRO:HD2	2.01	0.42
1:B:172:PRO:HG2	1:B:180:GLU:HG2	2.02	0.42
1:A:232[A]:LYS:HG3	1:A:459:PRO:HG2	2.02	0.42
1:A:172:PRO:HG2	1:A:180:GLU:HG2	2.02	0.42
1:A:206:ASN:HB3	1:A:310:GLN:HB2	2.02	0.42
1:B:16:HIS:HE1	5:B:563:HOH:O	2.02	0.42
1:A:232[B]:LYS:HD2	1:A:459:PRO:HG2	2.01	0.41
1:A:127:LEU:HD12	1:A:333:PHE:HB3	2.02	0.41
1:B:17:MET:O	1:B:24:VAL:HA	2.21	0.40
1:B:48:ASN:HD22	1:B:49:ARG:N	2.20	0.40
1:B:253[B]:ARG:CG	1:B:253[B]:ARG:NH1	2.41	0.40
1:A:143:GLN:HE21	1:A:291:VAL:HG21	1.86	0.40
1:B:24:VAL:HG21	1:B:49:ARG:HG2	2.04	0.40
1:B:67:LEU:HA	1:B:68:PRO:HD2	1.98	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 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	492/502 (98%)	482 (98%)	10 (2%)	0	100	100
1	B	500/502 (100%)	490 (98%)	10 (2%)	0	100	100
All	All	992/1004 (99%)	972 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	429/430 (100%)	416 (97%)	13 (3%)	41	41
1	B	437/430 (102%)	425 (97%)	12 (3%)	44	46
All	All	866/860 (101%)	841 (97%)	25 (3%)	43	43

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	40	ASN
1	A	48	ASN
1	A	82	ARG
1	A	93	ASN
1	A	129	TYR
1	A	185	ASN
1	A	309	ARG
1	A	317	TRP
1	A	324	LYS
1	A	332	LEU
1	A	395	TYR
1	A	484	LYS
1	B	11	LEU
1	B	40	ASN
1	B	48	ASN
1	B	93	ASN
1	B	94	LYS

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Mol	Chain	Res	Type
1	B	95	LEU
1	B	112	GLU
1	B	185	ASN
1	B	232[A]	LYS
1	B	232[B]	LYS
1	B	317	TRP
1	B	395	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	40	ASN
1	A	48	ASN
1	A	55	GLN
1	A	93	ASN
1	A	110	HIS
1	A	143	GLN
1	A	155	GLN
1	A	185	ASN
1	A	196	ASN
1	A	310	GLN
1	A	372	GLN
1	A	376	ASN
1	A	420	GLN
1	A	425	ASN
1	A	489	GLN
1	B	16	HIS
1	B	28	ASN
1	B	40	ASN
1	B	48	ASN
1	B	55	GLN
1	B	93	ASN
1	B	98	GLN
1	B	110	HIS
1	B	143	GLN
1	B	155	GLN
1	B	185	ASN
1	B	196	ASN
1	B	236	GLN
1	B	358	ASN
1	B	376	ASN

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Mol	Chain	Res	Type
1	B	420	GLN

### 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 ⓘ

4 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	LGU	C	1	2	13,13,13	1.07	0	18,19,19	0.65	0
2	BEM	C	2	2	12,12,13	0.92	0	14,17,19	0.62	0
2	LGU	D	1	2	13,13,13	1.12	0	18,19,19	0.60	0
2	BEM	D	2	2	12,12,13	0.79	0	14,17,19	0.65	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
2	LGU	C	1	2	-	1/4/24/24	0/1/1/1
2	BEM	C	2	2	-	0/4/21/24	0/1/1/1
2	LGU	D	1	2	-	1/4/24/24	0/1/1/1
2	BEM	D	2	2	-	0/4/21/24	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

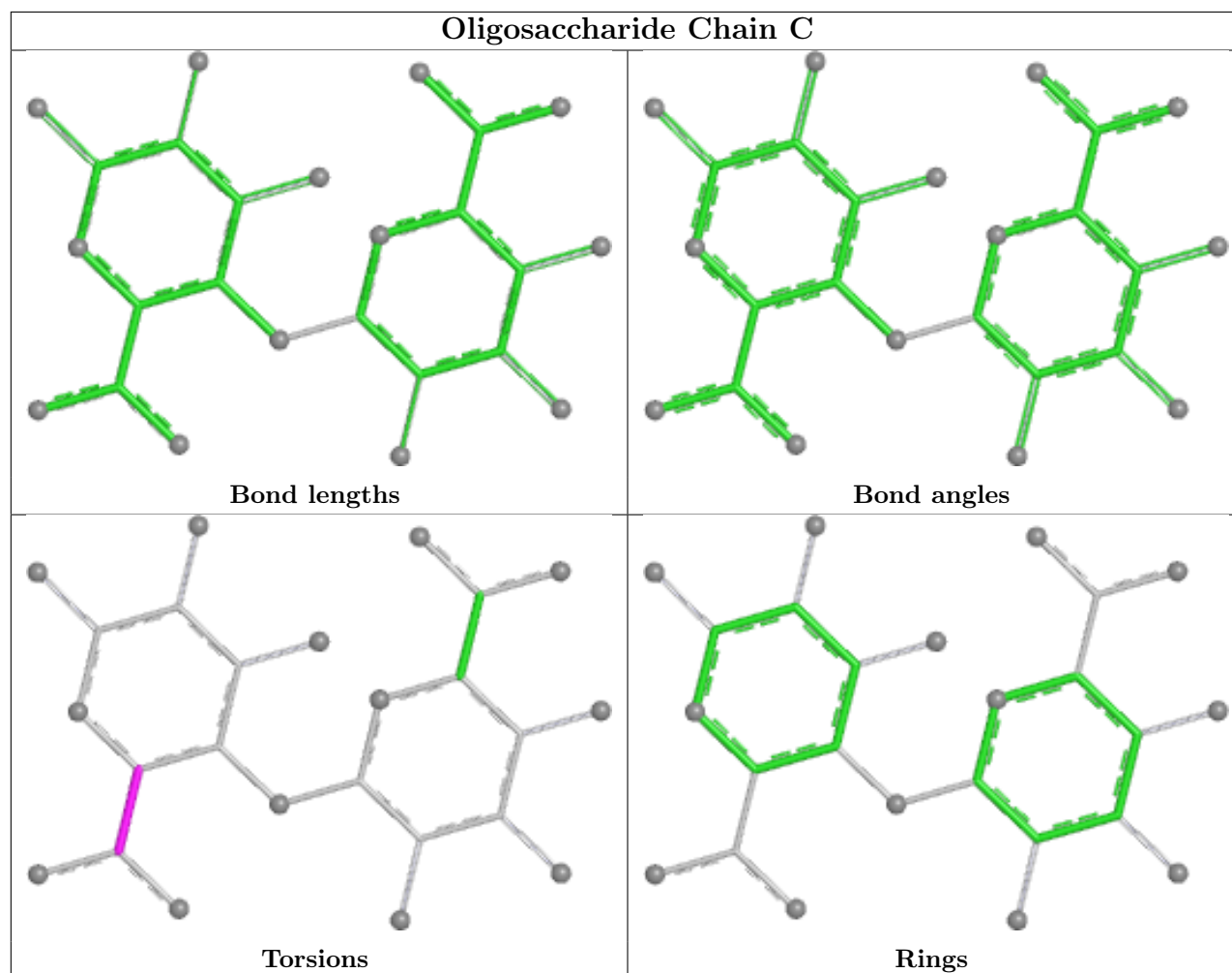
All (2) torsion outliers are listed below:

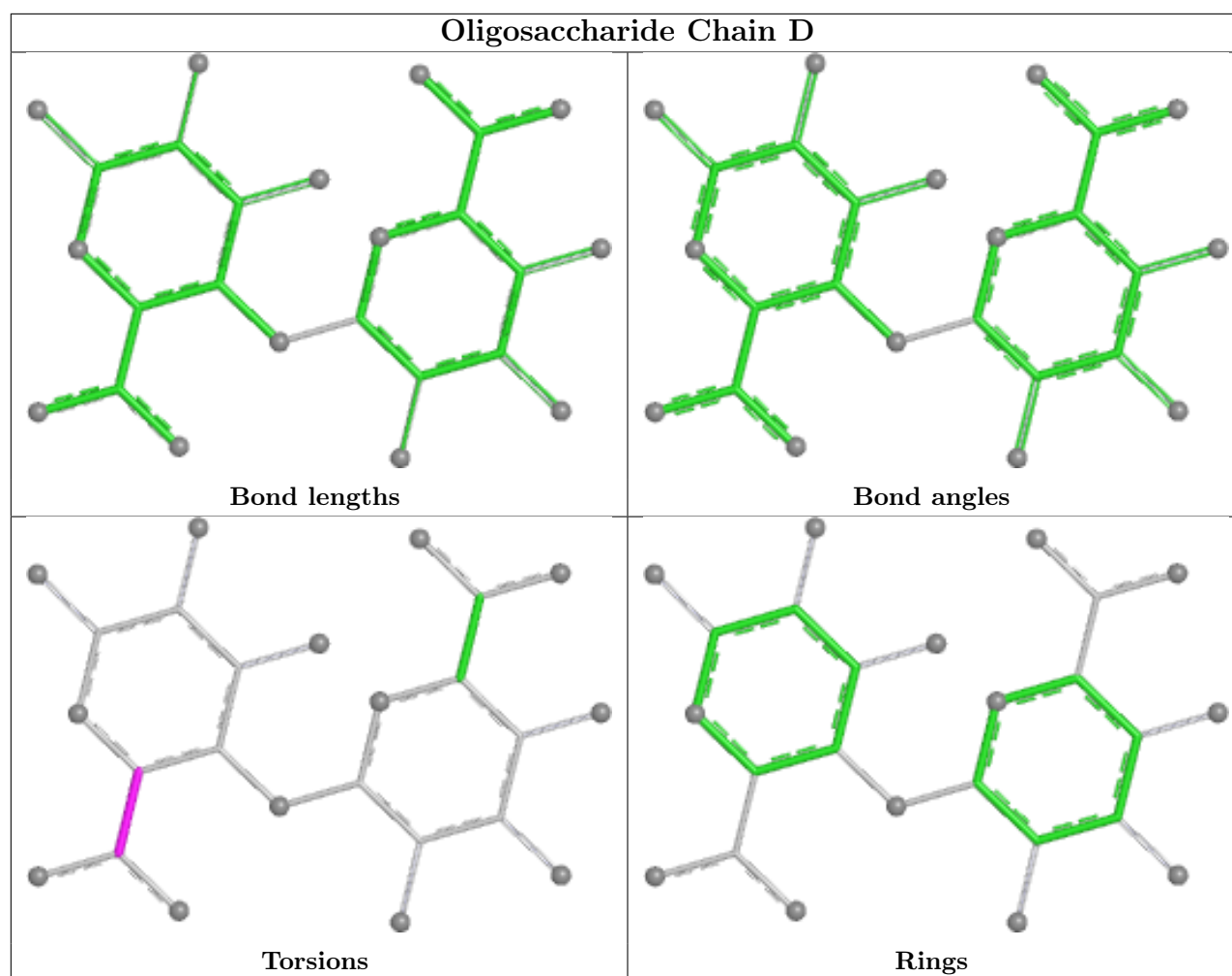
Mol	Chain	Res	Type	Atoms
2	C	1	LGU	O5-C5-C6-O6B
2	D	1	LGU	O5-C5-C6-O6B

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](#)

Of 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 for Mogul analysis.

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	$\# Z  > 2$
4	GOL	B	507	-	5,5,5	0.35	0	5,5,5	0.42	0
4	GOL	A	512	-	5,5,5	0.39	0	5,5,5	0.27	0
4	GOL	B	509	-	5,5,5	0.35	0	5,5,5	0.37	0
4	GOL	A	509	-	5,5,5	0.39	0	5,5,5	0.24	0
4	GOL	B	511	-	5,5,5	0.40	0	5,5,5	0.30	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	510	-	5,5,5	0.37	0	5,5,5	0.32	0
4	GOL	B	506	-	5,5,5	0.35	0	5,5,5	0.27	0
4	GOL	B	510	-	5,5,5	0.35	0	5,5,5	0.33	0
4	GOL	B	512	-	5,5,5	0.38	0	5,5,5	0.41	0
4	GOL	B	514	-	5,5,5	0.38	0	5,5,5	0.25	0
4	GOL	A	507	-	5,5,5	0.39	0	5,5,5	0.22	0
4	GOL	A	508	-	5,5,5	0.40	0	5,5,5	0.33	0
4	GOL	B	516	-	5,5,5	0.39	0	5,5,5	0.28	0
4	GOL	B	513	-	5,5,5	0.35	0	5,5,5	0.35	0
4	GOL	A	506	-	5,5,5	0.38	0	5,5,5	0.34	0
4	GOL	B	508	-	5,5,5	0.38	0	5,5,5	0.26	0
4	GOL	B	515	-	5,5,5	0.40	0	5,5,5	0.21	0
4	GOL	A	511	-	5,5,5	0.40	0	5,5,5	0.33	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
4	GOL	B	507	-	-	0/4/4/4	-
4	GOL	A	512	-	-	4/4/4/4	-
4	GOL	B	509	-	-	4/4/4/4	-
4	GOL	A	509	-	-	0/4/4/4	-
4	GOL	B	511	-	-	2/4/4/4	-
4	GOL	A	510	-	-	2/4/4/4	-
4	GOL	B	506	-	-	3/4/4/4	-
4	GOL	B	510	-	-	0/4/4/4	-
4	GOL	B	512	-	-	2/4/4/4	-
4	GOL	B	514	-	-	2/4/4/4	-
4	GOL	A	507	-	-	2/4/4/4	-
4	GOL	A	508	-	-	2/4/4/4	-
4	GOL	B	516	-	-	2/4/4/4	-
4	GOL	B	513	-	-	2/4/4/4	-
4	GOL	A	506	-	-	0/4/4/4	-
4	GOL	B	508	-	-	2/4/4/4	-
4	GOL	B	515	-	-	0/4/4/4	-
4	GOL	A	511	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	508	GOL	O1-C1-C2-O2
4	A	508	GOL	O1-C1-C2-C3
4	A	511	GOL	C1-C2-C3-O3
4	A	512	GOL	O1-C1-C2-O2
4	A	512	GOL	O1-C1-C2-C3
4	A	512	GOL	C1-C2-C3-O3
4	B	506	GOL	C1-C2-C3-O3
4	B	509	GOL	O1-C1-C2-C3
4	B	509	GOL	C1-C2-C3-O3
4	B	511	GOL	C1-C2-C3-O3
4	B	512	GOL	C1-C2-C3-O3
4	B	512	GOL	O2-C2-C3-O3
4	B	513	GOL	O1-C1-C2-C3
4	B	514	GOL	C1-C2-C3-O3
4	B	516	GOL	O1-C1-C2-C3
4	A	511	GOL	O2-C2-C3-O3
4	B	513	GOL	O1-C1-C2-O2
4	B	514	GOL	O2-C2-C3-O3
4	A	507	GOL	C1-C2-C3-O3
4	A	510	GOL	O1-C1-C2-C3
4	B	508	GOL	O1-C1-C2-C3
4	A	512	GOL	O2-C2-C3-O3
4	B	506	GOL	O2-C2-C3-O3
4	B	508	GOL	O1-C1-C2-O2
4	B	509	GOL	O1-C1-C2-O2
4	B	516	GOL	O1-C1-C2-O2
4	B	509	GOL	O2-C2-C3-O3
4	B	511	GOL	O2-C2-C3-O3
4	B	506	GOL	O1-C1-C2-C3
4	A	510	GOL	O1-C1-C2-O2
4	A	507	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	507	GOL	1	0
4	B	510	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	511	GOL	1	0

## 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	489/502 (97%)	-0.21	9 (1%) 68 66	15, 23, 34, 51	0
1	B	489/502 (97%)	-0.20	8 (1%) 72 70	16, 23, 32, 43	1 (0%)
All	All	978/1004 (97%)	-0.21	17 (1%) 70 68	15, 23, 33, 51	1 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	ALA	5.9
1	B	50	ALA	3.8
1	A	5	TRP	3.5
1	A	65	GLY	3.0
1	A	66	GLN	3.0
1	A	283	LYS	2.9
1	B	283	LYS	2.8
1	A	67	LEU	2.8
1	A	63	ALA	2.7
1	B	282	SER	2.6
1	B	371	ALA	2.5
1	A	3	ALA	2.5
1	B	8	GLU	2.5
1	B	66	GLN	2.2
1	B	490	TYR	2.2
1	A	490	TYR	2.0
1	B	353	THR	2.0

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

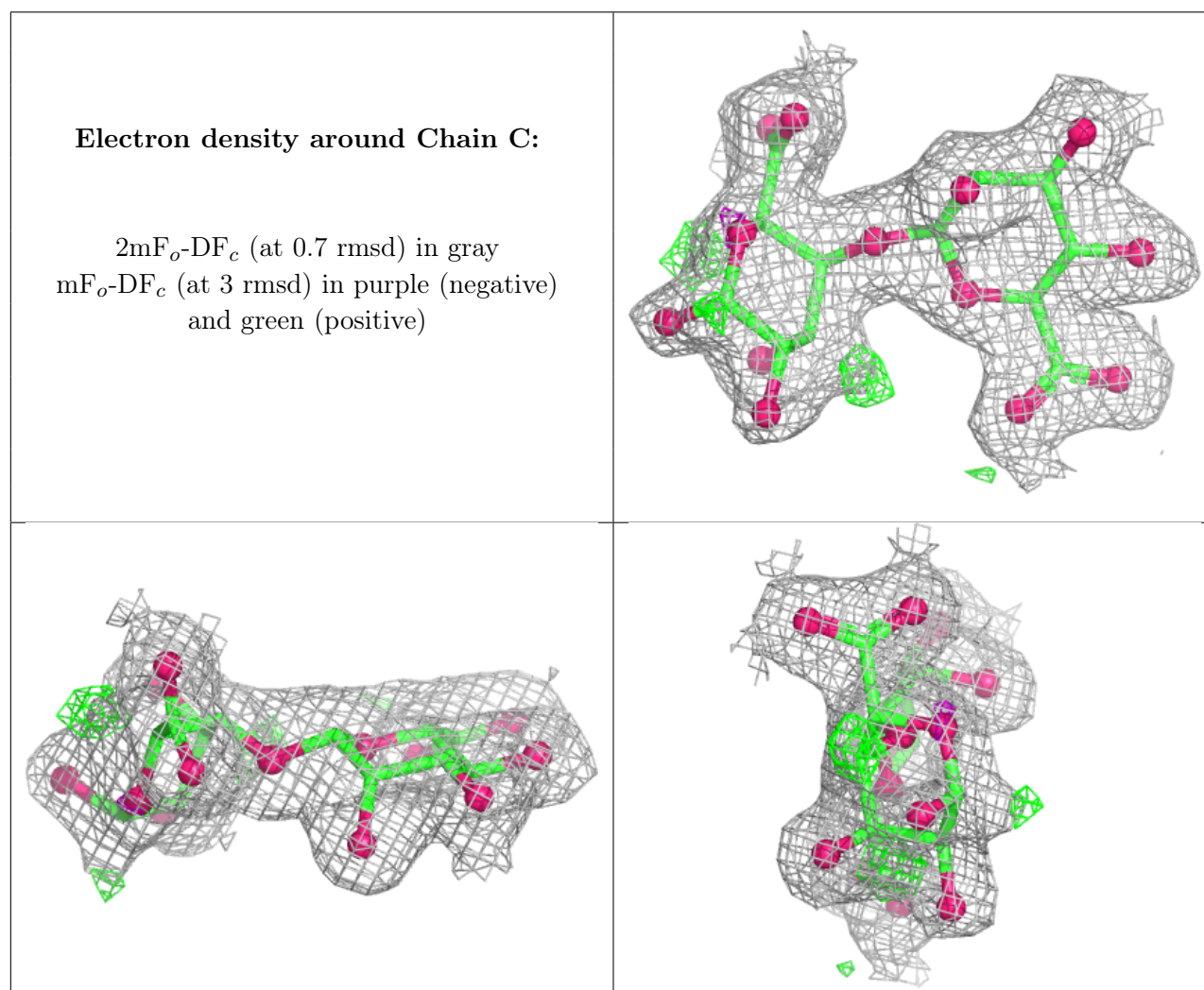
There are no non-standard protein/DNA/RNA residues in this entry.

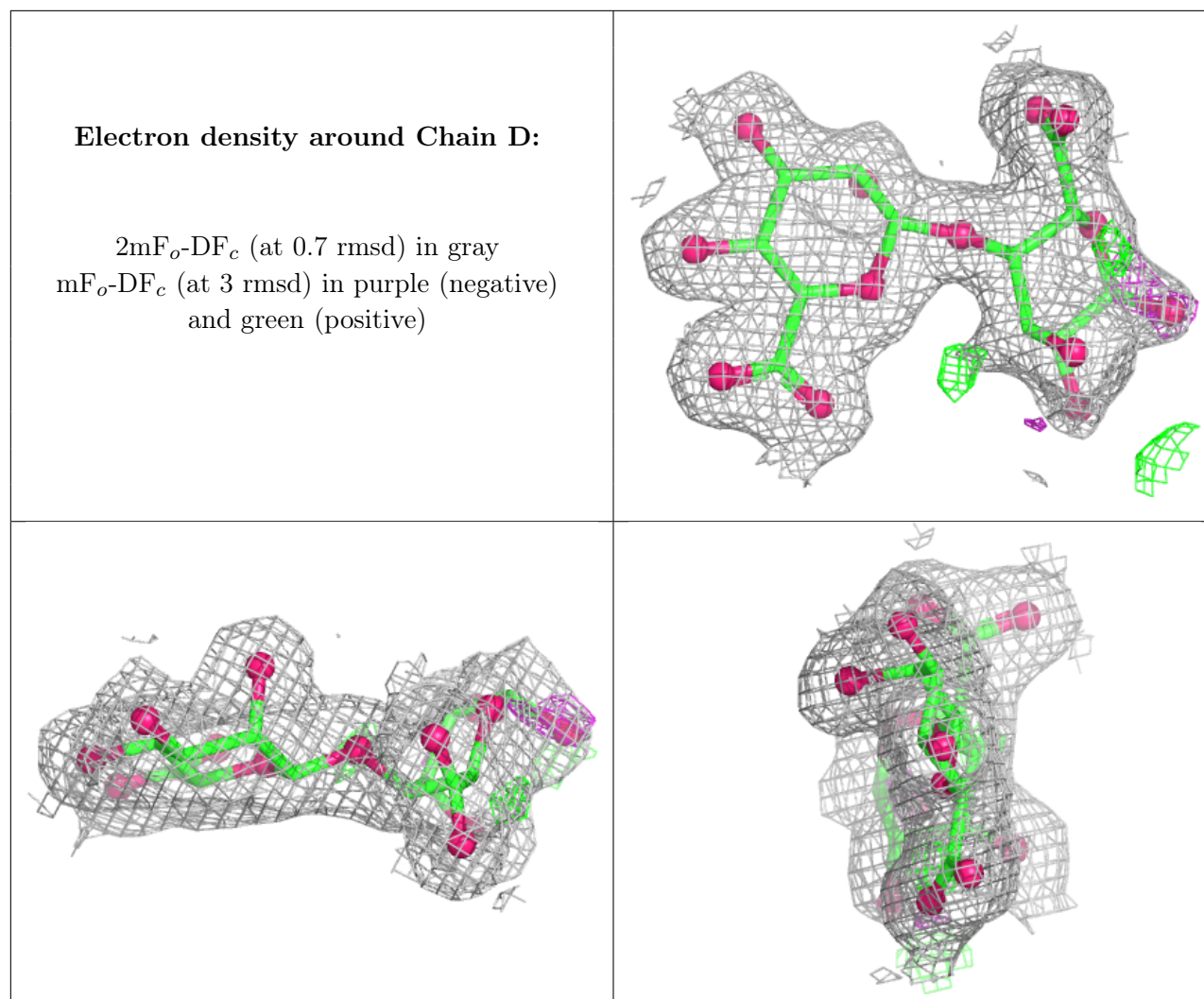
### 6.3 Carbohydrates [i](#)

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( $\text{\AA}^2$ )	Q<0.9
2	LGU	D	1	13/13	0.86	0.14	21,23,26,27	0
2	LGU	C	1	13/13	0.88	0.14	23,25,28,29	0
2	BEM	C	2	12/13	0.97	0.09	18,19,20,20	0
2	BEM	D	2	12/13	0.97	0.10	17,17,18,18	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 ⓘ

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( $\text{\AA}^2$ )	Q<0.9
4	GOL	B	516	6/6	0.65	0.43	64,65,65,65	0
4	GOL	A	512	6/6	0.67	0.30	62,62,62,62	0
4	GOL	B	511	6/6	0.70	0.35	59,59,59,59	0
4	GOL	B	507	6/6	0.72	0.20	38,40,40,40	0
4	GOL	B	510	6/6	0.76	0.22	53,53,53,53	0
4	GOL	B	508	6/6	0.77	0.20	56,56,56,57	0
4	GOL	A	507	6/6	0.77	0.23	51,52,52,52	0
4	GOL	B	513	6/6	0.79	0.26	42,43,43,44	0
4	GOL	B	506	6/6	0.81	0.21	36,36,36,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	GOL	B	515	6/6	0.82	0.18	30,31,31,31	0
4	GOL	A	508	6/6	0.82	0.21	37,39,39,40	0
4	GOL	B	514	6/6	0.83	0.19	52,53,53,53	0
4	GOL	B	512	6/6	0.83	0.15	38,39,39,39	0
4	GOL	B	509	6/6	0.83	0.18	47,49,49,49	0
4	GOL	A	510	6/6	0.84	0.19	36,37,38,38	0
4	GOL	A	506	6/6	0.86	0.20	32,33,33,33	0
4	GOL	A	511	6/6	0.88	0.15	43,44,44,45	0
4	GOL	A	509	6/6	0.96	0.09	27,27,27,28	0
3	CA	B	503	1/1	0.98	0.08	24,24,24,24	0
3	CA	A	503	1/1	0.99	0.06	22,22,22,22	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.