



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 30, 2023 – 04:23 PM JST

PDB ID : 4U5K  
Title : Complex structure of mutant CtCel5E (E314A) with cellobiose  
Authors : Guo, R.T.; Huang, C.H.; Wu, T.H.  
Deposited on : 2014-07-25  
Resolution : 2.65 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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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

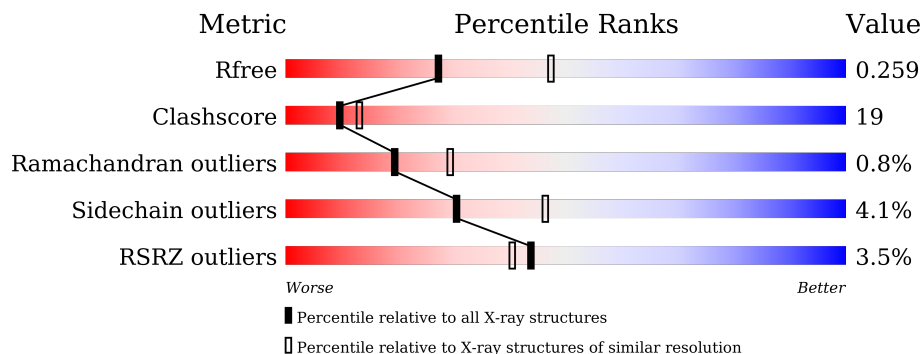
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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	403	
1	B	403	
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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BGC	C	2	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5575 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 Endoglucanase H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	304	2532	1625	424	472	11	0	0	0
1	B	298	2473	1588	411	463	11	0	0	0

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P16218
A	2	GLY	-	expression tag	UNP P16218
A	3	SER	-	expression tag	UNP P16218
A	4	SER	-	expression tag	UNP P16218
A	5	HIS	-	expression tag	UNP P16218
A	6	HIS	-	expression tag	UNP P16218
A	7	HIS	-	expression tag	UNP P16218
A	8	HIS	-	expression tag	UNP P16218
A	9	HIS	-	expression tag	UNP P16218
A	10	HIS	-	expression tag	UNP P16218
A	11	SER	-	expression tag	UNP P16218
A	12	SER	-	expression tag	UNP P16218
A	13	GLY	-	expression tag	UNP P16218
A	14	LEU	-	expression tag	UNP P16218
A	15	VAL	-	expression tag	UNP P16218
A	16	PRO	-	expression tag	UNP P16218
A	17	ARG	-	expression tag	UNP P16218
A	18	GLY	-	expression tag	UNP P16218
A	19	SER	-	expression tag	UNP P16218
A	20	HIS	-	expression tag	UNP P16218
A	21	MET	-	expression tag	UNP P16218
A	22	ALA	-	expression tag	UNP P16218
A	23	SER	-	expression tag	UNP P16218
A	24	MET	-	expression tag	UNP P16218
A	25	THR	-	expression tag	UNP P16218

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Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	expression tag	UNP P16218
A	27	GLY	-	expression tag	UNP P16218
A	28	GLN	-	expression tag	UNP P16218
A	29	GLN	-	expression tag	UNP P16218
A	30	MET	-	expression tag	UNP P16218
A	31	GLY	-	expression tag	UNP P16218
A	32	ARG	-	expression tag	UNP P16218
A	33	ILE	-	expression tag	UNP P16218
A	34	GLU	-	expression tag	UNP P16218
A	35	GLY	-	expression tag	UNP P16218
A	36	ARG	-	expression tag	UNP P16218
A	37	GLU	-	expression tag	UNP P16218
A	38	PHE	-	expression tag	UNP P16218
A	314	ALA	GLU	engineered mutation	UNP P16218
B	1	MET	-	expression tag	UNP P16218
B	2	GLY	-	expression tag	UNP P16218
B	3	SER	-	expression tag	UNP P16218
B	4	SER	-	expression tag	UNP P16218
B	5	HIS	-	expression tag	UNP P16218
B	6	HIS	-	expression tag	UNP P16218
B	7	HIS	-	expression tag	UNP P16218
B	8	HIS	-	expression tag	UNP P16218
B	9	HIS	-	expression tag	UNP P16218
B	10	HIS	-	expression tag	UNP P16218
B	11	SER	-	expression tag	UNP P16218
B	12	SER	-	expression tag	UNP P16218
B	13	GLY	-	expression tag	UNP P16218
B	14	LEU	-	expression tag	UNP P16218
B	15	VAL	-	expression tag	UNP P16218
B	16	PRO	-	expression tag	UNP P16218
B	17	ARG	-	expression tag	UNP P16218
B	18	GLY	-	expression tag	UNP P16218
B	19	SER	-	expression tag	UNP P16218
B	20	HIS	-	expression tag	UNP P16218
B	21	MET	-	expression tag	UNP P16218
B	22	ALA	-	expression tag	UNP P16218
B	23	SER	-	expression tag	UNP P16218
B	24	MET	-	expression tag	UNP P16218
B	25	THR	-	expression tag	UNP P16218
B	26	GLY	-	expression tag	UNP P16218
B	27	GLY	-	expression tag	UNP P16218
B	28	GLN	-	expression tag	UNP P16218

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Chain	Residue	Modelled	Actual	Comment	Reference
B	29	GLN	-	expression tag	UNP P16218
B	30	MET	-	expression tag	UNP P16218
B	31	GLY	-	expression tag	UNP P16218
B	32	ARG	-	expression tag	UNP P16218
B	33	ILE	-	expression tag	UNP P16218
B	34	GLU	-	expression tag	UNP P16218
B	35	GLY	-	expression tag	UNP P16218
B	36	ARG	-	expression tag	UNP P16218
B	37	GLU	-	expression tag	UNP P16218
B	38	PHE	-	expression tag	UNP P16218
B	314	ALA	GLU	engineered mutation	UNP P16218

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



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

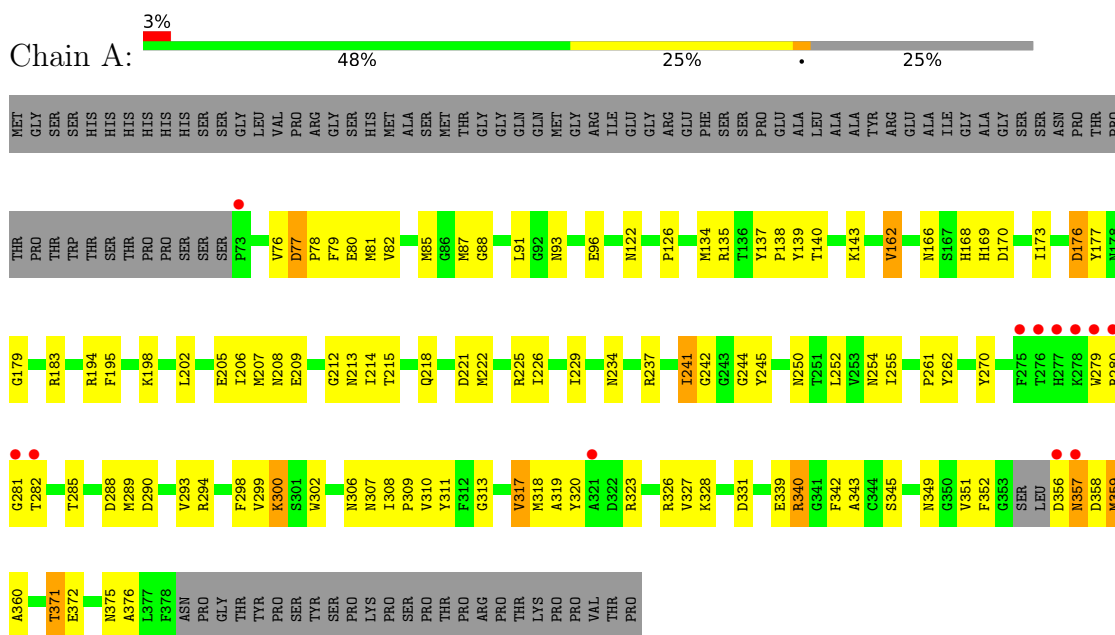
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	284	Total	O	0	0
			284	284		
3	B	240	Total	O	0	0
			240	240		

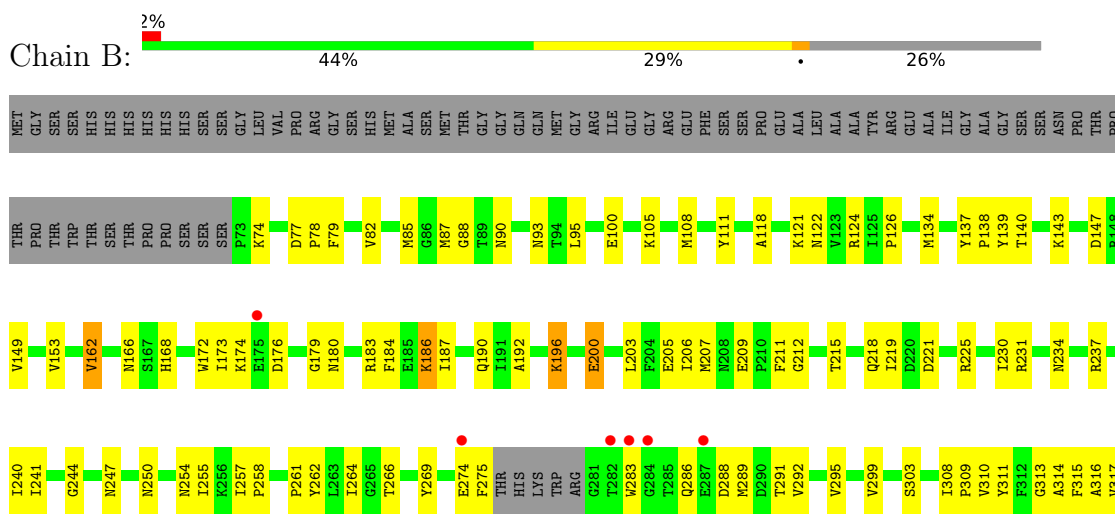
### 3 Residue-property plots [i](#)

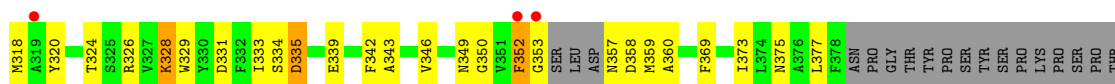
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Endoglucanase H



#### • Molecule 1: Endoglucanase H





PRO  
ARG  
PRO  
THR  
LYS  
PRO  
PRO  
VAL  
THR  
PRO

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

Chain C: 100%

BGC1  
BGC2

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

Chain D: 50% 50%

BGC1  
BGC2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.94Å 74.94Å 254.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.65 24.98 – 2.65	Depositor EDS
% Data completeness (in resolution range)	94.0 (25.00-2.65) 94.2 (24.98-2.65)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.24 (at 2.64Å)	Xtrriage
Refinement program	CNS 1.21	Depositor
R, $R_{free}$	0.183 , 0.255 0.185 , 0.259	Depositor DCC
$R_{free}$ test set	1000 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.1	Xtrriage
Anisotropy	0.080	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 57.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5575	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: BGC

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	0.38	0/2605	0.64	2/3533 (0.1%)
1	B	0.35	0/2542	0.60	1/3446 (0.0%)
All	All	0.37	0/5147	0.62	3/6979 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	162	VAL	N-CA-C	-5.70	95.61	111.00
1	A	176	ASP	N-CA-C	-5.38	96.48	111.00
1	A	162	VAL	N-CA-C	-5.20	96.96	111.00

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	2532	0	2392	93	0
1	B	2473	0	2337	103	0
2	C	23	0	20	2	0
2	D	23	0	20	2	0
3	A	284	0	0	7	0
3	B	240	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5575	0	4769	190	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 19.

All (190) 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:82:VAL:HG21	1:A:309:PRO:HB2	1.26	1.17
1:B:134:MET:HE2	1:B:140:THR:H	1.14	1.07
1:B:82:VAL:HG21	1:B:309:PRO:HB2	1.41	1.01
1:B:352:PHE:HA	1:B:360:ALA:HB3	1.54	0.88
1:B:315:PHE:CZ	1:B:333:ILE:HG22	2.10	0.86
1:A:255:ILE:HG21	1:B:291:THR:HG23	1.59	0.85
1:A:85:MET:HG2	1:A:162:VAL:HG11	1.60	0.83
1:A:170:ASP:O	1:A:173:ILE:HG22	1.82	0.80
1:B:315:PHE:HB3	1:B:346:VAL:HA	1.67	0.76
1:A:82:VAL:HG21	1:A:309:PRO:CB	2.14	0.75
1:A:134:MET:HE2	1:A:140:THR:HB	1.68	0.75
1:A:318:MET:HG2	3:A:770:HOH:O	1.86	0.73
1:A:206:ILE:HD12	1:A:241:ILE:HG22	1.69	0.73
1:A:285:THR:HG22	1:A:288:ASP:OD2	1.89	0.73
1:B:134:MET:HE2	1:B:140:THR:N	1.98	0.73
1:A:285:THR:HG23	1:A:288:ASP:H	1.56	0.71
1:A:356:ASP:O	1:A:357:ASN:HB2	1.91	0.70
1:A:349:ASN:HD22	1:A:351:VAL:HB	1.56	0.70
1:B:357:ASN:HD22	1:B:359:MET:HG3	1.56	0.69
1:A:87:MET:HE2	1:A:88:GLY:N	2.08	0.69
1:B:349:ASN:HD21	1:B:359:MET:HB2	1.58	0.68
1:A:134:MET:HE2	1:A:140:THR:CB	2.23	0.68
1:A:135:ARG:HG3	3:A:608:HOH:O	1.93	0.68
1:A:349:ASN:ND2	1:A:351:VAL:HB	2.08	0.68
1:A:280:ARG:HH22	1:A:288:ASP:CG	1.97	0.68
1:B:134:MET:CE	1:B:140:THR:H	2.01	0.68
1:B:105:LYS:HD2	1:B:111:TYR:OH	1.94	0.67
1:B:291:THR:O	1:B:295:VAL:HG23	1.93	0.67
1:A:166:ASN:HB3	1:A:205:GLU:HB3	1.75	0.67
1:B:317:VAL:HG21	2:D:1:BGC:H6C2	1.77	0.67
1:B:82:VAL:HG21	1:B:309:PRO:CB	2.20	0.66
1:A:215:THR:OG1	1:A:218:GLN:HG3	1.95	0.66
1:A:208:ASN:ND2	1:A:209:GLU:HG3	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:MET:HE2	1:A:140:THR:H	1.63	0.64
1:B:286:GLN:HA	1:B:289:MET:HG2	1.79	0.64
1:B:82:VAL:HG13	1:B:343:ALA:HB2	1.79	0.64
1:A:214:ILE:HG13	1:A:218:GLN:OE1	1.98	0.63
1:B:257:ILE:HD13	1:B:308:ILE:HD12	1.81	0.63
1:A:134:MET:CE	1:A:140:THR:H	2.13	0.61
1:B:326:ARG:NH1	1:B:373:ILE:HD11	2.16	0.61
1:B:90:ASN:ND2	1:B:313:GLY:O	2.34	0.60
1:B:87:MET:HE2	1:B:88:GLY:O	2.01	0.60
1:B:317:VAL:HG22	1:B:317:VAL:O	2.02	0.60
1:B:85:MET:HG2	1:B:162:VAL:HG11	1.84	0.59
1:A:313:GLY:O	1:A:345:SER:HB2	2.02	0.59
1:B:138:PRO:O	1:B:186:LYS:HG2	2.02	0.59
1:A:331:ASP:HB2	1:A:376:ALA:HB1	1.85	0.59
1:B:329:TRP:O	1:B:333:ILE:HG12	2.02	0.59
1:A:87:MET:HE2	1:A:88:GLY:H	1.68	0.59
1:A:198:LYS:O	1:A:237:ARG:NH2	2.36	0.59
1:A:351:VAL:H	1:A:360:ALA:HB2	1.68	0.59
1:A:241:ILE:H	1:A:241:ILE:HD13	1.67	0.58
1:A:176:ASP:OD2	1:A:179:GLY:HA3	2.03	0.58
1:A:255:ILE:HD13	1:B:295:VAL:HG22	1.86	0.58
1:A:307:ASN:O	1:A:308:ILE:HD13	2.04	0.58
1:A:143:LYS:HB2	3:A:676:HOH:O	2.04	0.58
1:A:169:HIS:HE1	3:A:873:HOH:O	1.86	0.58
1:A:82:VAL:HG13	1:A:343:ALA:HB2	1.85	0.57
1:A:252:LEU:C	1:A:252:LEU:HD23	2.24	0.57
1:B:139:TYR:HB2	1:B:186:LYS:HG2	1.84	0.57
1:B:186:LYS:HA	1:B:186:LYS:HE3	1.85	0.57
1:B:310:VAL:HB	1:B:342:PHE:CD1	2.39	0.57
1:B:85:MET:O	1:B:121:LYS:HB3	2.03	0.57
1:B:95:LEU:HD23	1:B:100:GLU:HG3	1.87	0.56
1:B:174:LYS:HZ2	1:B:212:GLY:HA3	1.70	0.56
1:A:198:LYS:HG2	1:A:202:LEU:HD22	1.87	0.55
1:A:285:THR:HG22	1:A:288:ASP:CG	2.27	0.55
1:A:317:VAL:HG22	2:C:1:BGC:H6C2	1.88	0.55
1:B:315:PHE:HZ	1:B:333:ILE:HG22	1.69	0.55
1:A:280:ARG:NH2	1:A:288:ASP:OD2	2.38	0.55
1:A:96:GLU:HB2	1:A:168:HIS:HB3	1.90	0.54
1:B:206:ILE:HD12	1:B:241:ILE:HG22	1.89	0.54
1:B:274:GLU:O	1:B:275:PHE:C	2.45	0.54
1:B:221:ASP:O	1:B:225:ARG:HG2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:ILE:HG12	1:B:309:PRO:HG2	1.89	0.54
1:A:208:ASN:CG	1:A:209:GLU:HG3	2.28	0.54
1:B:350:GLY:HA2	3:B:700:HOH:O	2.06	0.53
1:A:300:LYS:HD3	1:A:340:ARG:HG2	1.91	0.53
1:A:137:TYR:CD1	1:A:138:PRO:HA	2.44	0.52
1:B:318:MET:HB2	1:B:320:TYR:CZ	2.44	0.52
1:A:331:ASP:HA	1:A:376:ALA:O	2.10	0.52
1:B:139:TYR:CD1	1:B:183:ARG:HG3	2.44	0.52
1:A:212:GLY:O	1:A:213:ASN:HB2	2.10	0.51
1:B:358:ASP:O	1:B:359:MET:HB2	2.10	0.51
1:A:173:ILE:HG21	1:A:207:MET:HE2	1.91	0.51
1:B:172:TRP:CZ3	1:B:180:ASN:HB3	2.46	0.51
1:B:326:ARG:HH12	1:B:373:ILE:HD11	1.74	0.51
1:B:172:TRP:HZ3	1:B:180:ASN:HB3	1.76	0.51
1:B:196:LYS:HA	1:B:234:ASN:OD1	2.11	0.51
1:A:310:VAL:HB	1:A:342:PHE:CD1	2.46	0.50
1:B:93:ASN:HD22	1:B:93:ASN:N	2.09	0.50
1:B:74:LYS:O	1:B:200:GLU:HG2	2.11	0.50
1:B:317:VAL:CG2	2:D:1:BGC:H6C2	2.42	0.50
1:B:314:ALA:HB2	3:B:719:HOH:O	2.11	0.49
1:B:234:ASN:OD1	1:B:237:ARG:NH1	2.45	0.49
1:A:311:TYR:HA	1:A:343:ALA:O	2.13	0.49
1:B:88:GLY:HA3	1:B:122:ASN:CG	2.33	0.49
1:B:137:TYR:CG	1:B:138:PRO:HA	2.48	0.49
1:B:303:SER:HB2	1:B:310:VAL:HG23	1.94	0.49
1:A:261:PRO:HG2	1:A:262:TYR:CD2	2.48	0.49
1:A:352:PHE:HA	1:A:357:ASN:HA	1.93	0.49
1:B:215:THR:OG1	1:B:218:GLN:HG3	2.13	0.49
1:A:299:VAL:HG12	1:A:310:VAL:HG21	1.95	0.48
1:A:282:THR:HB	1:A:319:ALA:O	2.13	0.48
1:B:283:TRP:CD1	1:B:318:MET:HB3	2.48	0.48
1:A:244:GLY:HA3	1:A:252:LEU:O	2.13	0.48
1:A:173:ILE:O	1:A:176:ASP:O	2.32	0.48
1:A:255:ILE:HG13	1:B:250:ASN:HA	1.96	0.48
1:A:234:ASN:OD1	1:A:237:ARG:NH1	2.45	0.48
1:A:281:GLY:O	1:A:282:THR:OG1	2.31	0.47
1:B:215:THR:O	1:B:219:ILE:HG13	2.14	0.47
1:B:231:ARG:NH2	1:B:262:TYR:HB2	2.29	0.47
1:A:317:VAL:O	1:A:317:VAL:HG13	2.13	0.47
1:A:302:TRP:CD1	1:A:306:ASN:ND2	2.83	0.47
1:B:320:TYR:HD2	1:B:326:ARG:HD2	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:ASP:O	1:A:225:ARG:HG2	2.14	0.47
1:B:328:LYS:HE3	1:B:328:LYS:HA	1.97	0.47
1:B:318:MET:HB2	1:B:320:TYR:CE1	2.50	0.47
1:A:96:GLU:HB3	1:A:126:PRO:HB2	1.97	0.46
1:B:299:VAL:HG12	1:B:310:VAL:HG21	1.97	0.46
1:B:78:PRO:O	1:B:82:VAL:HG23	2.14	0.46
1:B:357:ASN:ND2	1:B:359:MET:HG3	2.27	0.46
1:B:211:PHE:CZ	1:B:244:GLY:HA2	2.50	0.46
1:A:139:TYR:CD1	1:A:183:ARG:HG3	2.51	0.46
1:A:250:ASN:HA	1:B:255:ILE:HG12	1.98	0.46
1:B:118:ALA:HB2	1:B:369:PHE:CZ	2.50	0.46
1:A:357:ASN:HB3	3:A:817:HOH:O	2.16	0.46
1:B:149:VAL:O	1:B:153:VAL:HG23	2.15	0.46
1:B:173:ILE:HD12	1:B:184:PHE:HB2	1.98	0.46
1:B:266:THR:HA	1:B:311:TYR:O	2.16	0.45
1:A:255:ILE:CG2	1:B:291:THR:HG23	2.38	0.45
1:B:88:GLY:HA3	1:B:122:ASN:OD1	2.16	0.45
1:B:288:ASP:O	1:B:292:VAL:HG23	2.16	0.45
1:B:205:GLU:HA	1:B:240:ILE:HB	1.98	0.45
1:A:320:TYR:CD2	1:A:326:ARG:HG3	2.52	0.44
1:A:323:ARG:O	1:A:327:VAL:HG23	2.17	0.44
1:B:183:ARG:O	1:B:187:ILE:HG13	2.16	0.44
1:B:353:GLY:HA2	1:B:358:ASP:CB	2.47	0.44
1:B:331:ASP:O	1:B:335:ASP:HB2	2.17	0.44
1:A:222:MET:O	1:A:226:ILE:HG12	2.18	0.44
1:A:294:ARG:HG3	1:A:298:PHE:CE2	2.53	0.44
1:B:230:ILE:O	1:B:234:ASN:HB2	2.18	0.44
1:A:93:ASN:HD21	2:C:2:BGC:H2	1.82	0.43
1:A:166:ASN:OD1	1:A:166:ASN:C	2.57	0.43
1:B:334:SER:OG	1:B:377:LEU:HD23	2.17	0.43
1:A:371:THR:HG22	1:A:372:GLU:N	2.34	0.43
1:A:137:TYR:CG	1:A:138:PRO:HA	2.53	0.43
1:A:242:GLY:O	1:A:254:ASN:ND2	2.51	0.43
1:B:139:TYR:CD2	1:B:183:ARG:HA	2.54	0.43
1:B:93:ASN:OD1	1:B:349:ASN:HB2	2.19	0.43
1:B:335:ASP:O	1:B:339:GLU:HB2	2.19	0.43
1:A:77:ASP:OD2	1:A:79:PHE:HB2	2.19	0.43
1:B:143:LYS:O	1:B:147:ASP:OD1	2.36	0.42
1:A:290:ASP:O	1:A:293:VAL:HG22	2.18	0.42
1:A:255:ILE:CD1	1:B:295:VAL:HG22	2.49	0.42
1:A:289:MET:O	1:A:293:VAL:HG13	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:PRO:HG2	3:A:653:HOH:O	2.20	0.42
1:A:225:ARG:O	1:A:229:ILE:HG13	2.19	0.42
1:B:349:ASN:ND2	1:B:359:MET:HB2	2.29	0.42
1:A:78:PRO:O	1:A:82:VAL:HG23	2.20	0.42
1:B:93:ASN:HD22	1:B:93:ASN:H	1.67	0.42
1:B:93:ASN:N	1:B:93:ASN:ND2	2.68	0.42
1:A:371:THR:HB	3:A:777:HOH:O	2.20	0.42
1:B:166:ASN:HB3	1:B:205:GLU:HB3	2.01	0.42
1:B:124:ARG:O	1:B:126:PRO:HD3	2.19	0.41
1:B:261:PRO:HG2	1:B:262:TYR:CD2	2.55	0.41
1:A:88:GLY:HA3	1:A:122:ASN:CG	2.40	0.41
1:A:299:VAL:CG1	1:A:310:VAL:HG21	2.51	0.41
1:A:76:VAL:HG13	1:A:80:GLU:CD	2.41	0.41
1:B:95:LEU:CD2	1:B:100:GLU:HG3	2.51	0.41
1:B:328:LYS:HE3	1:B:328:LYS:CA	2.49	0.41
1:A:195:PHE:O	1:A:198:LYS:HB2	2.20	0.41
1:A:245:TYR:HB2	1:A:270:TYR:HB3	2.02	0.41
1:B:134:MET:CE	1:B:138:PRO:HD2	2.50	0.41
1:B:257:ILE:HA	1:B:258:PRO:HD3	1.96	0.41
1:A:245:TYR:CG	1:A:270:TYR:HB3	2.56	0.41
1:A:359:MET:H	1:A:359:MET:HG2	1.77	0.41
1:B:77:ASP:OD2	1:B:79:PHE:HB2	2.21	0.41
1:B:192:ALA:HB1	1:B:234:ASN:HD22	1.86	0.41
1:B:247:ASN:HA	1:B:269:TYR:CE2	2.56	0.41
1:B:324:THR:O	1:B:328:LYS:HB2	2.21	0.41
1:A:91:LEU:HD12	1:A:91:LEU:HA	1.93	0.40
1:B:139:TYR:O	1:B:190:GLN:NE2	2.51	0.40
1:B:315:PHE:CZ	1:B:333:ILE:CG2	2.95	0.40
1:A:81:MET:O	1:A:85:MET:HG3	2.21	0.40
1:A:323:ARG:NH1	1:A:372:GLU:HB2	2.37	0.40
1:B:126:PRO:HB3	1:B:168:HIS:ND1	2.37	0.40
1:B:134:MET:HE1	1:B:138:PRO:HD2	2.03	0.40
1:B:173:ILE:HG21	1:B:207:MET:CE	2.51	0.40
1:B:176:ASP:OD2	1:B:179:GLY:HA3	2.21	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	300/403 (74%)	275 (92%)	22 (7%)	3 (1%)	15	23
1	B	292/403 (72%)	266 (91%)	24 (8%)	2 (1%)	22	33
All	All	592/806 (73%)	541 (91%)	46 (8%)	5 (1%)	19	29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	ASN
1	A	358	ASP
1	A	177	TYR
1	B	196	LYS
1	B	316	ALA

### 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	270/353 (76%)	258 (96%)	12 (4%)	28	43
1	B	264/353 (75%)	254 (96%)	10 (4%)	33	49
All	All	534/706 (76%)	512 (96%)	22 (4%)	30	46

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	ASP

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Mol	Chain	Res	Type
1	A	194	ARG
1	A	241	ILE
1	A	279	TRP
1	A	300	LYS
1	A	317	VAL
1	A	328	LYS
1	A	339	GLU
1	A	340	ARG
1	A	359	MET
1	A	371	THR
1	A	375	ASN
1	B	108	MET
1	B	186	LYS
1	B	200	GLU
1	B	203	LEU
1	B	209	GLU
1	B	254	ASN
1	B	328	LYS
1	B	335	ASP
1	B	352	PHE
1	B	375	ASN

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

Mol	Chain	Res	Type
1	A	93	ASN
1	A	169	HIS
1	A	180	ASN
1	A	277	HIS
1	A	349	ASN
1	B	93	ASN
1	B	180	ASN
1	B	286	GLN
1	B	357	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](#)

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

## 5.5 Carbohydrates [i](#)

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	BGC	C	1	2	12,12,12	3.07	3 (25%)	17,17,17	1.17	1 (5%)
2	BGC	C	2	2	11,11,12	1.72	4 (36%)	15,15,17	0.85	0
2	BGC	D	1	2	12,12,12	3.14	3 (25%)	17,17,17	1.16	1 (5%)
2	BGC	D	2	2	11,11,12	1.60	2 (18%)	15,15,17	0.94	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	BGC	C	1	2	-	2/2/22/22	0/1/1/1
2	BGC	C	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1
2	BGC	D	2	2	-	0/2/19/22	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	BGC	O6-C6	-9.43	1.02	1.42
2	C	1	BGC	O6-C6	-9.42	1.02	1.42
2	D	2	BGC	O5-C1	3.55	1.49	1.43
2	D	1	BGC	C4-C5	3.40	1.60	1.53
2	C	2	BGC	O5-C1	3.09	1.48	1.43
2	C	1	BGC	C4-C5	2.96	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	BGC	O5-C5	2.49	1.48	1.43
2	C	2	BGC	O5-C5	2.33	1.48	1.43
2	C	2	BGC	C2-C3	2.33	1.55	1.52
2	D	1	BGC	O5-C1	2.26	1.48	1.42
2	C	2	BGC	C4-C5	2.20	1.57	1.53
2	C	1	BGC	O5-C1	2.15	1.48	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	BGC	O6-C6-C5	3.20	122.28	111.29
2	D	1	BGC	O6-C6-C5	3.12	122.00	111.29

There are no chirality outliers.

All (2) torsion outliers are listed below:

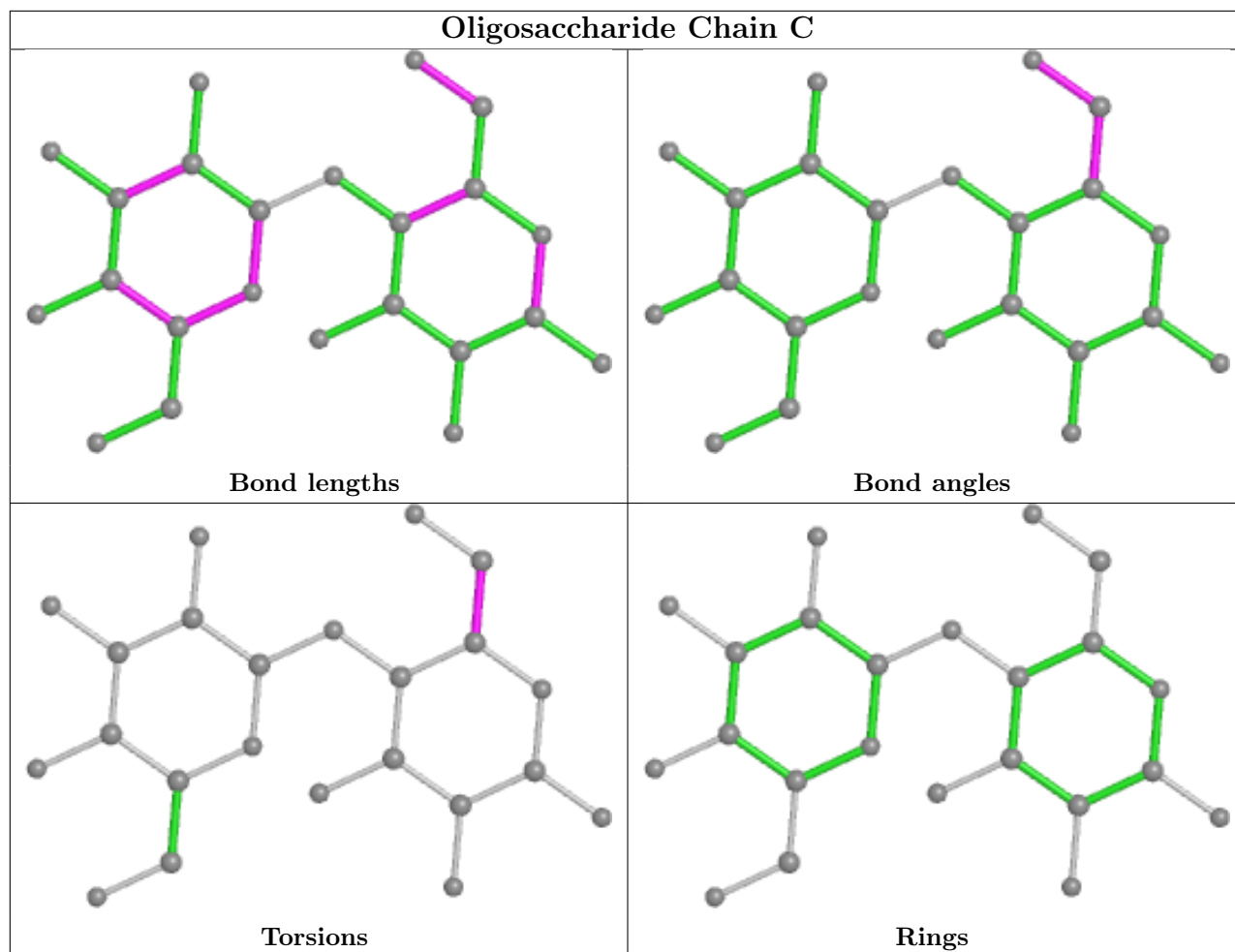
Mol	Chain	Res	Type	Atoms
2	C	1	BGC	C4-C5-C6-O6
2	C	1	BGC	O5-C5-C6-O6

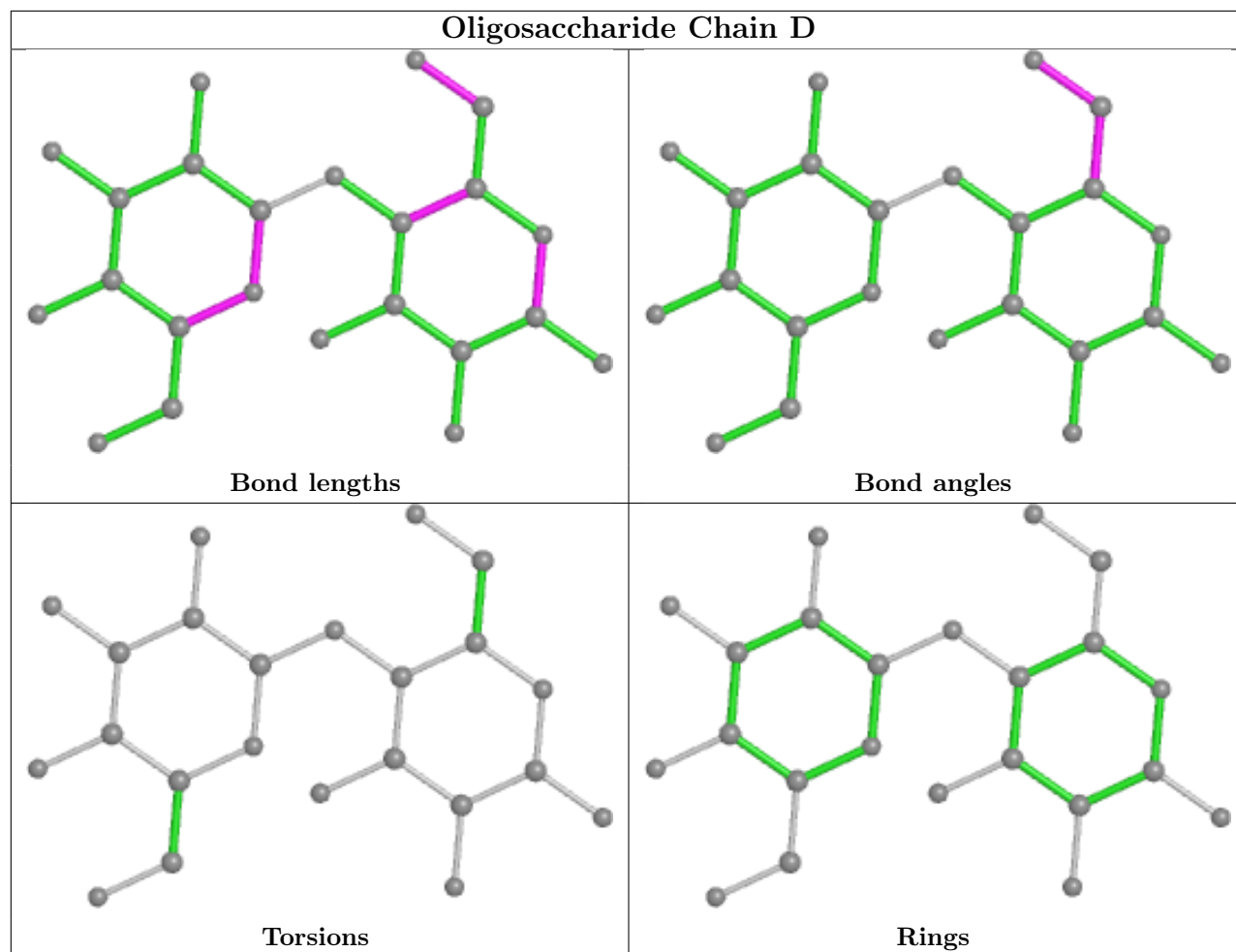
There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	BGC	1	0
2	D	1	BGC	2	0
2	C	2	BGC	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

### 6.1 Protein, DNA and RNA chains

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	304/403 (75%)	-0.36	12 (3%) 39 35	16, 32, 64, 103	0
1	B	298/403 (73%)	-0.19	9 (3%) 50 47	22, 40, 78, 100	0
All	All	602/806 (74%)	-0.28	21 (3%) 44 40	16, 35, 74, 103	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	357	ASN	6.1
1	A	73	PRO	4.8
1	A	279	TRP	4.8
1	A	275	PHE	3.9
1	A	280	ARG	3.7
1	A	356	ASP	3.6
1	A	321	ALA	3.6
1	A	281	GLY	3.2
1	B	282	THR	3.2
1	B	353	GLY	2.9
1	B	284	GLY	2.8
1	A	282	THR	2.7
1	B	283	TRP	2.6
1	A	276	THR	2.3
1	B	319	ALA	2.3
1	B	274	GLU	2.3
1	B	175	GLU	2.3
1	B	352	PHE	2.1
1	A	277	HIS	2.0
1	A	278	LYS	2.0
1	B	287	GLU	2.0

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

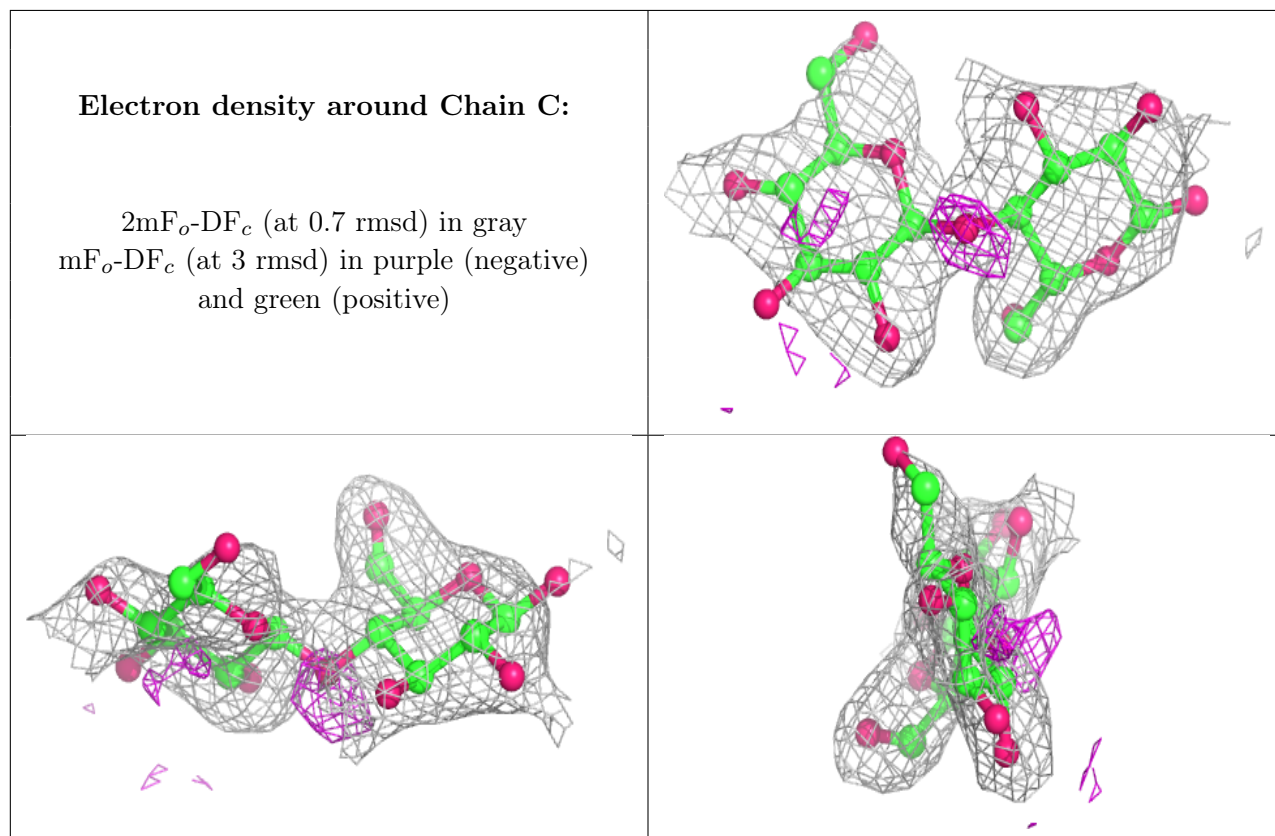
There are no non-standard protein/DNA/RNA residues in this entry.

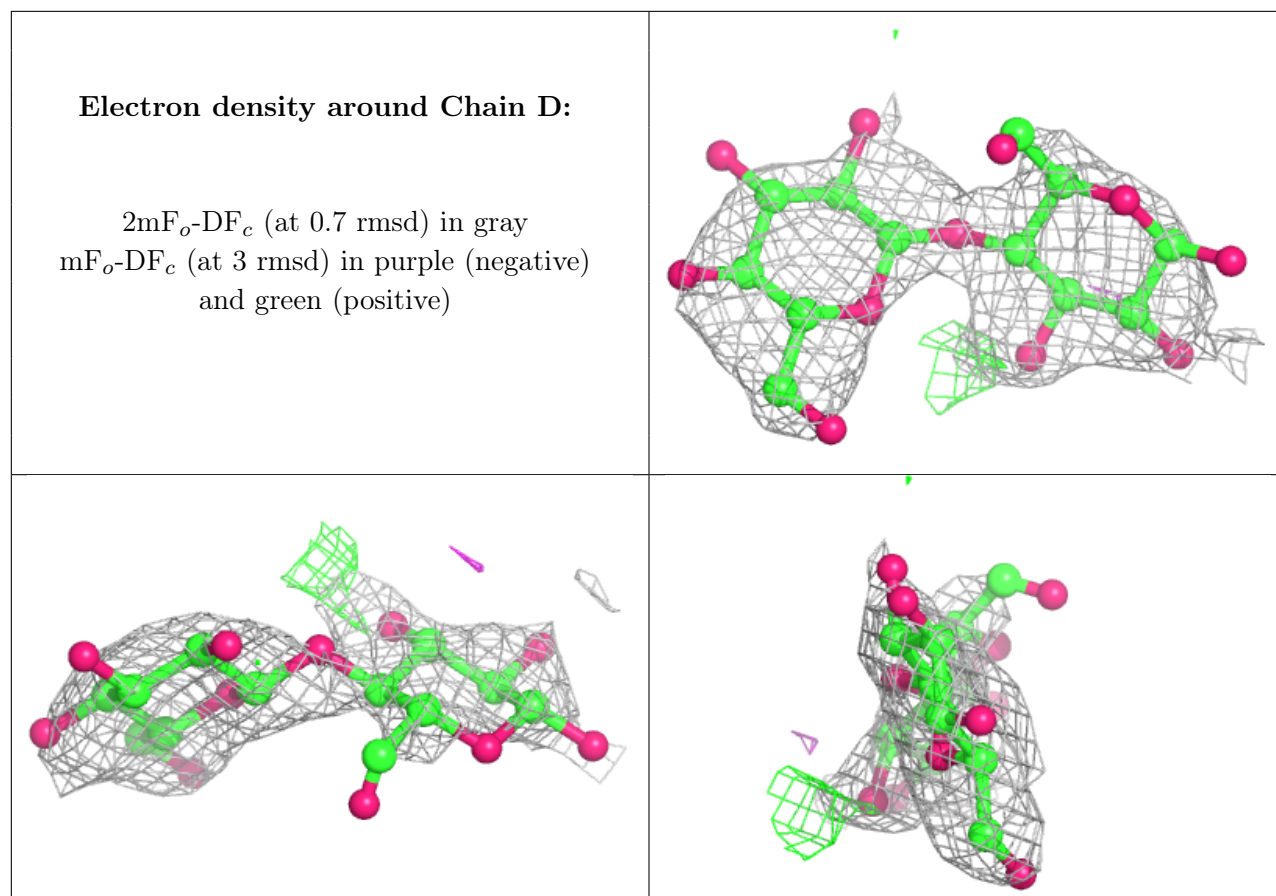
## 6.3 Carbohydrates [i](#)

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(Å <sup>2</sup> )	Q<0.9
2	BGC	D	1	12/12	0.78	0.35	100,101,102,102	0
2	BGC	C	2	11/12	0.79	0.41	64,68,70,70	0
2	BGC	D	2	11/12	0.81	0.29	96,98,98,99	0
2	BGC	C	1	12/12	0.87	0.21	57,62,65,66	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 [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.