



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 8, 2024 – 06:36 am GMT

PDB ID : 5NZ8  
Title : Clostridium thermocellum cellodextrin phosphorylase with cellotetraose and phosphate bound  
Authors : O'Neill, E.C.; Pergolizzi, G.; Stevenson, C.E.M.; Lawson, D.M.; Nepogodiev, S.A.; Field, R.A.  
Deposited on : 2017-05-12  
Resolution : 3.00 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 : 1.8.4, 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

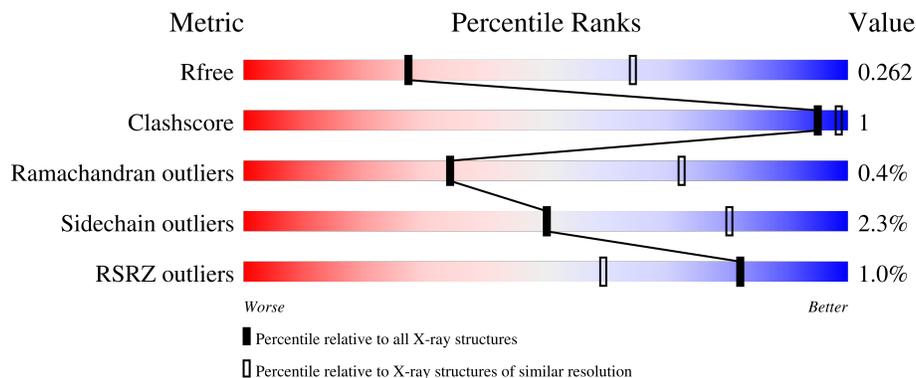
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

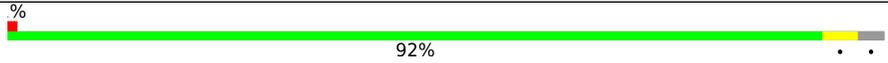
The reported resolution of this entry is 3.00 Å.

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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	1009	 83% 12%
1	B	1009	 92%
2	C	4	 50% 50%
2	D	4	 75% 25%

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	BGC	D	1	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 14127 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 Cellodextrin phosphorylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	884	6696	4316	1116	1237	27	0	0	0
1	B	974	7331	4693	1225	1382	31	0	0	0

There are 50 discrepancies between the modelled and reference sequences:

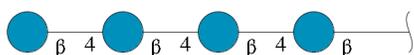
Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP Q93HT8
A	-23	GLY	-	expression tag	UNP Q93HT8
A	-22	SER	-	expression tag	UNP Q93HT8
A	-21	SER	-	expression tag	UNP Q93HT8
A	-20	HIS	-	expression tag	UNP Q93HT8
A	-19	HIS	-	expression tag	UNP Q93HT8
A	-18	HIS	-	expression tag	UNP Q93HT8
A	-17	HIS	-	expression tag	UNP Q93HT8
A	-16	HIS	-	expression tag	UNP Q93HT8
A	-15	HIS	-	expression tag	UNP Q93HT8
A	-14	SER	-	expression tag	UNP Q93HT8
A	-13	SER	-	expression tag	UNP Q93HT8
A	-12	GLY	-	expression tag	UNP Q93HT8
A	-11	LEU	-	expression tag	UNP Q93HT8
A	-10	VAL	-	expression tag	UNP Q93HT8
A	-9	PRO	-	expression tag	UNP Q93HT8
A	-8	ARG	-	expression tag	UNP Q93HT8
A	-7	GLY	-	expression tag	UNP Q93HT8
A	-6	SER	-	expression tag	UNP Q93HT8
A	-5	HIS	-	expression tag	UNP Q93HT8
A	-4	MET	-	expression tag	UNP Q93HT8
A	-3	LEU	-	expression tag	UNP Q93HT8
A	-2	GLU	-	expression tag	UNP Q93HT8
A	-1	ASP	-	expression tag	UNP Q93HT8
A	0	PRO	-	expression tag	UNP Q93HT8

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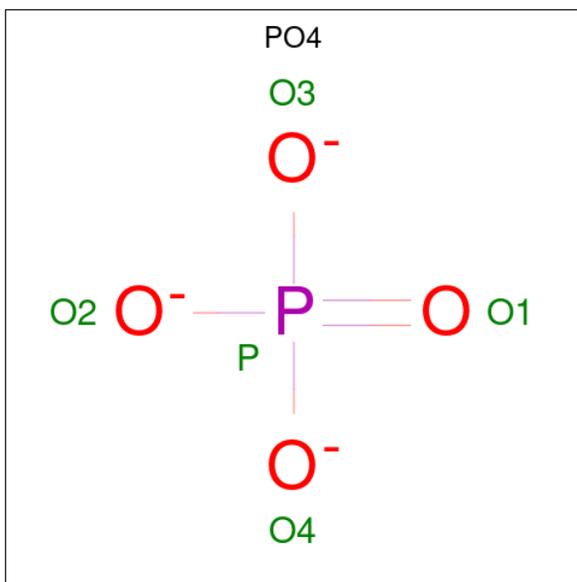
Chain	Residue	Modelled	Actual	Comment	Reference
B	-24	MET	-	initiating methionine	UNP Q93HT8
B	-23	GLY	-	expression tag	UNP Q93HT8
B	-22	SER	-	expression tag	UNP Q93HT8
B	-21	SER	-	expression tag	UNP Q93HT8
B	-20	HIS	-	expression tag	UNP Q93HT8
B	-19	HIS	-	expression tag	UNP Q93HT8
B	-18	HIS	-	expression tag	UNP Q93HT8
B	-17	HIS	-	expression tag	UNP Q93HT8
B	-16	HIS	-	expression tag	UNP Q93HT8
B	-15	HIS	-	expression tag	UNP Q93HT8
B	-14	SER	-	expression tag	UNP Q93HT8
B	-13	SER	-	expression tag	UNP Q93HT8
B	-12	GLY	-	expression tag	UNP Q93HT8
B	-11	LEU	-	expression tag	UNP Q93HT8
B	-10	VAL	-	expression tag	UNP Q93HT8
B	-9	PRO	-	expression tag	UNP Q93HT8
B	-8	ARG	-	expression tag	UNP Q93HT8
B	-7	GLY	-	expression tag	UNP Q93HT8
B	-6	SER	-	expression tag	UNP Q93HT8
B	-5	HIS	-	expression tag	UNP Q93HT8
B	-4	MET	-	expression tag	UNP Q93HT8
B	-3	LEU	-	expression tag	UNP Q93HT8
B	-2	GLU	-	expression tag	UNP Q93HT8
B	-1	ASP	-	expression tag	UNP Q93HT8
B	0	PRO	-	expression tag	UNP Q93HT8

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	0	0	0
			45	24	21			
2	D	4	Total	C	O	0	0	0
			45	24	21			

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

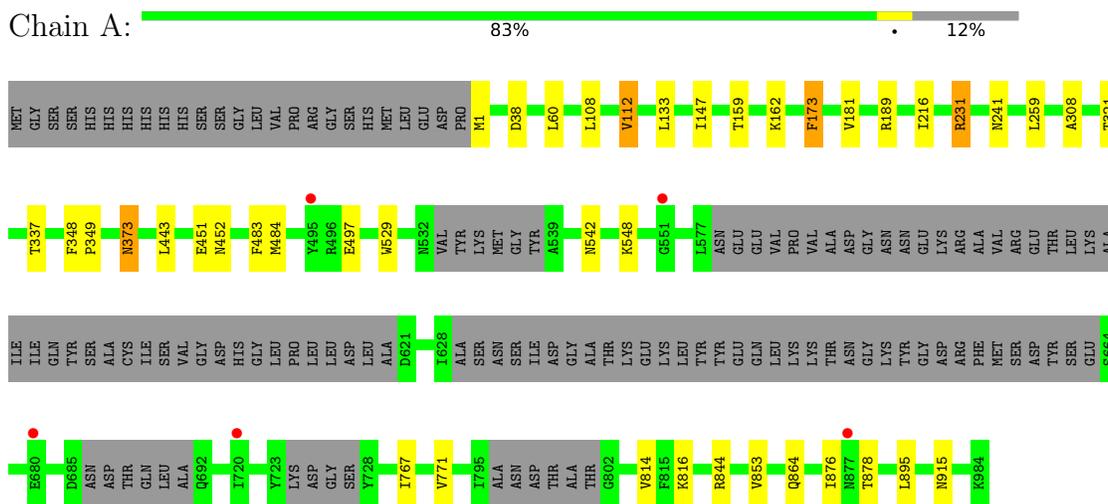


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

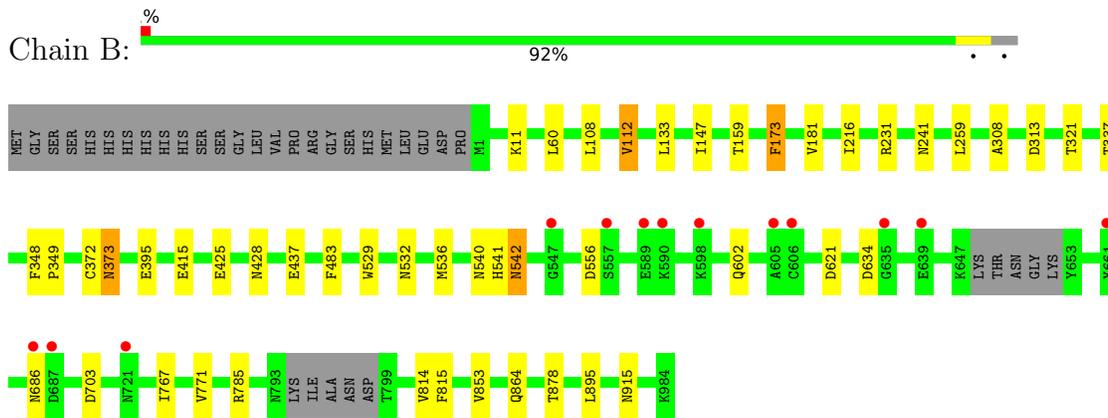
### 3 Residue-property plots [i](#)

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: Cellodextrin phosphorylase



- Molecule 1: Cellodextrin phosphorylase



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



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

Chain D:  75% 25%

BGC1  
BGC2  
BGC3  
BGC4

## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.81Å 152.97Å 92.22Å 90.00° 114.42° 90.00°	Depositor
Resolution (Å)	27.62 – 3.00 27.62 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.4 (27.62-3.00) 96.6 (27.62-3.00)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.57 (at 3.00Å)	Xtrriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.210 , 0.262 0.209 , 0.262	Depositor DCC
$R_{free}$ test set	2133 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	98.9	Xtrriage
Anisotropy	0.474	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 67.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.029 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14127	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	114.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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, PO4

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	0.41	0/6850	0.60	1/9326 (0.0%)
1	B	0.41	0/7497	0.59	1/10216 (0.0%)
All	All	0.41	0/14347	0.60	2/19542 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	621	ASP	CB-CG-OD1	6.04	123.74	118.30
1	A	189	ARG	NE-CZ-NH2	-5.14	117.73	120.30

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	6696	0	6187	16	0
1	B	7331	0	6728	16	0
2	C	45	0	39	0	0
2	D	45	0	39	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
All	All	14127	0	12993	32	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All (32) 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:541:HIS:CD2	1:B:542:ASN:HD21	2.14	0.64
1:B:373:ASN:HD22	1:B:373:ASN:N	2.03	0.56
1:A:373:ASN:HD22	1:A:373:ASN:N	2.04	0.56
1:B:173:PHE:CE2	1:B:181:VAL:HG11	2.46	0.51
1:B:60:LEU:HD21	1:B:108:LEU:HD23	1.93	0.49
1:B:308:ALA:HB3	1:B:321:THR:HG23	1.94	0.49
1:A:173:PHE:CE2	1:A:181:VAL:HG11	2.48	0.49
1:A:308:ALA:HB3	1:A:321:THR:HG23	1.95	0.48
1:A:60:LEU:HD21	1:A:108:LEU:HD23	1.96	0.47
1:B:112:VAL:O	1:B:112:VAL:HG12	2.15	0.47
1:A:112:VAL:HG12	1:A:112:VAL:O	2.15	0.47
1:A:162:LYS:HD3	1:A:484:MET:SD	2.55	0.46
1:A:159:THR:HB	1:A:259:LEU:HD11	1.97	0.46
1:B:532:ASN:HB3	1:B:540:ASN:HD22	1.81	0.46
1:B:159:THR:HB	1:B:259:LEU:HD11	1.98	0.46
1:A:814:VAL:HG23	1:A:878:THR:HG22	1.99	0.45
1:A:133:LEU:HD21	1:A:216:ILE:HG21	1.98	0.45
1:B:541:HIS:ND1	1:B:556:ASP:OD2	2.50	0.44
1:B:814:VAL:HG23	1:B:878:THR:HG22	1.99	0.44
1:B:133:LEU:HD21	1:B:216:ILE:HG21	2.00	0.43
1:B:348:PHE:HB3	1:B:349:PRO:HD2	2.00	0.43
1:A:231:ARG:HG2	1:A:443:LEU:HD11	2.01	0.42
1:B:147:ILE:HD11	1:B:483:PHE:CD2	2.54	0.42
1:A:767:ILE:O	1:A:771:VAL:HG23	2.20	0.42
1:B:853:VAL:HG12	1:B:895:LEU:HD22	2.02	0.41
1:A:147:ILE:HD11	1:A:483:PHE:CD2	2.55	0.41
1:A:348:PHE:HB3	1:A:349:PRO:HD2	2.03	0.41
1:A:816:LYS:HD2	1:A:876:ILE:HG13	2.03	0.41
1:A:542:ASN:HB2	1:A:548:LYS:O	2.21	0.41
1:A:853:VAL:HG12	1:A:895:LEU:HD22	2.03	0.40
1:B:425:GLU:HA	1:B:428:ASN:HD22	1.87	0.40
1:B:767:ILE:O	1:B:771:VAL:HG23	2.20	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	870/1009 (86%)	821 (94%)	46 (5%)	3 (0%)	41	76
1	B	968/1009 (96%)	909 (94%)	55 (6%)	4 (0%)	34	72
All	All	1838/2018 (91%)	1730 (94%)	101 (6%)	7 (0%)	34	72

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	11	LYS
1	A	241	ASN
1	B	241	ASN
1	A	497	GLU
1	B	686	ASN
1	A	112	VAL
1	B	112	VAL

### 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	646/872 (74%)	634 (98%)	12 (2%)	57	84
1	B	709/872 (81%)	690 (97%)	19 (3%)	44	77
All	All	1355/1744 (78%)	1324 (98%)	31 (2%)	50	80

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	38	ASP
1	A	173	PHE
1	A	231	ARG
1	A	337	THR
1	A	373	ASN
1	A	451	GLU
1	A	452	ASN
1	A	529	TRP
1	A	844	ARG
1	A	864	GLN
1	A	915	ASN
1	B	173	PHE
1	B	231	ARG
1	B	313	ASP
1	B	337	THR
1	B	372	CYS
1	B	373	ASN
1	B	395	GLU
1	B	415	GLU
1	B	437	GLU
1	B	529	TRP
1	B	536	MET
1	B	542	ASN
1	B	602	GLN
1	B	634	ASP
1	B	703	ASP
1	B	785	ARG
1	B	815	PHE
1	B	864	GLN
1	B	915	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	GLN
1	A	302	ASN
1	A	416	ASN
1	A	532	ASN
1	A	562	GLN
1	A	915	ASN
1	B	9	ASN
1	B	96	GLN

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Mol	Chain	Res	Type
1	B	190	GLN
1	B	302	ASN
1	B	357	ASN
1	B	428	ASN
1	B	532	ASN
1	B	540	ASN
1	B	542	ASN
1	B	562	GLN
1	B	915	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](#)

8 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	0.17	0	17,17,17	0.16	0
2	BGC	C	2	2	11,11,12	0.22	0	15,15,17	0.67	1 (6%)
2	BGC	C	3	2	11,11,12	0.21	0	15,15,17	0.44	0
2	BGC	C	4	2	11,11,12	0.41	0	15,15,17	1.21	2 (13%)
2	BGC	D	1	2	12,12,12	0.14	0	17,17,17	0.27	0
2	BGC	D	2	2	11,11,12	0.26	0	15,15,17	0.48	0
2	BGC	D	3	2	11,11,12	0.24	0	15,15,17	0.50	0
2	BGC	D	4	2	11,11,12	0.39	0	15,15,17	0.98	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	BGC	C	1	2	-	0/2/22/22	0/1/1/1
2	BGC	C	2	2	-	2/2/19/22	0/1/1/1
2	BGC	C	3	2	-	0/2/19/22	0/1/1/1
2	BGC	C	4	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
2	BGC	D	3	2	-	2/2/19/22	0/1/1/1
2	BGC	D	4	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	C	4	BGC	C1-O5-C5	2.86	116.07	112.19
2	D	4	BGC	C1-C2-C3	2.80	113.11	109.67
2	C	4	BGC	C1-C2-C3	2.77	113.08	109.67
2	C	2	BGC	C1-O5-C5	2.23	115.22	112.19

There are no chirality outliers.

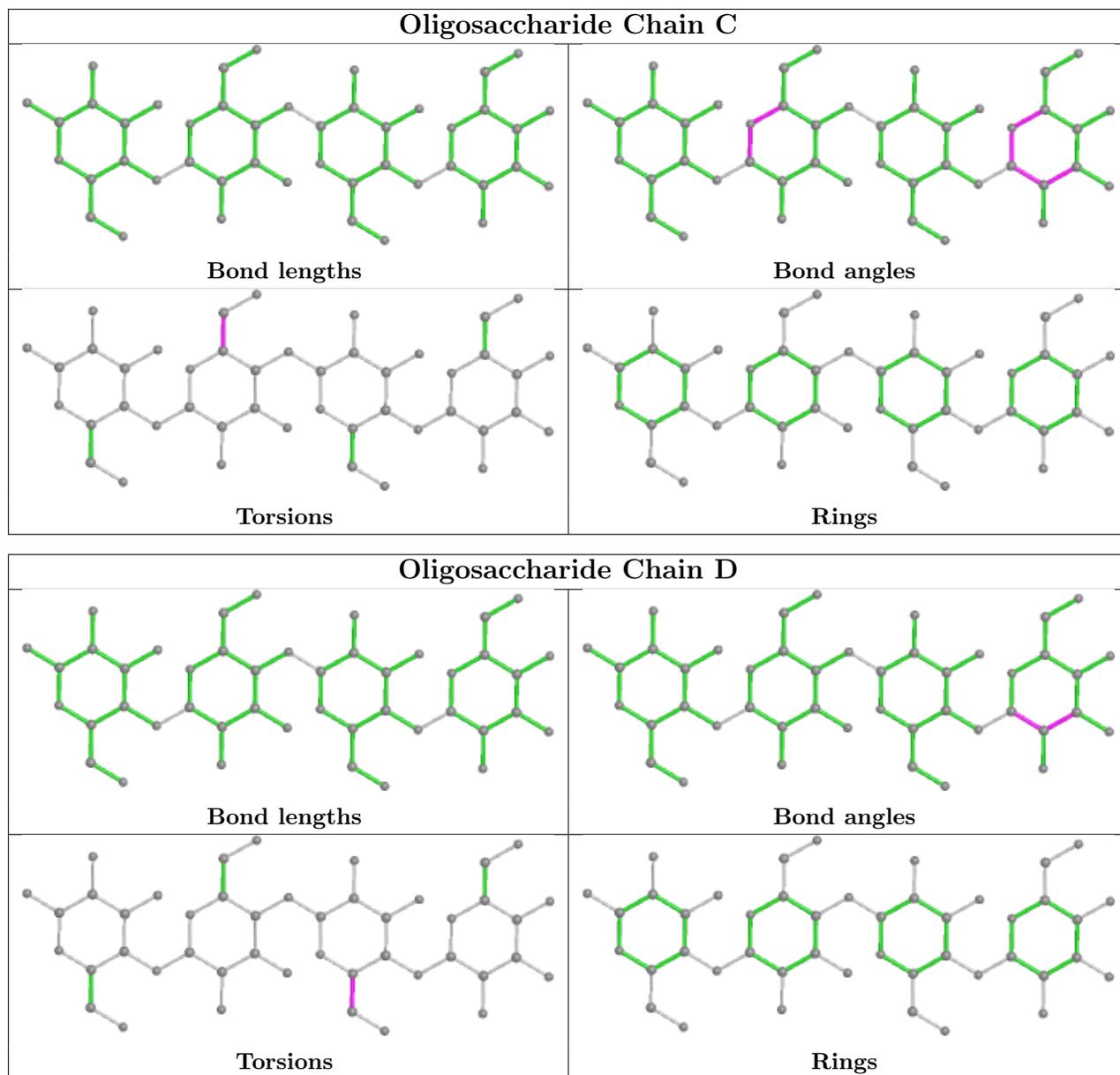
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	BGC	O5-C5-C6-O6
2	D	3	BGC	C4-C5-C6-O6
2	D	3	BGC	O5-C5-C6-O6
2	C	2	BGC	C4-C5-C6-O6

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	PO4	B	1005	-	4,4,4	1.35	1 (25%)	6,6,6	0.46	0
3	PO4	A	1005	-	4,4,4	1.31	1 (25%)	6,6,6	0.53	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1005	PO4	P-O1	2.59	1.56	1.50
3	A	1005	PO4	P-O1	2.52	1.56	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 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	884/1009 (87%)	-0.29	5 (0%) 89 72	74, 110, 153, 197	0
1	B	974/1009 (96%)	-0.25	13 (1%) 77 51	70, 114, 161, 208	0
All	All	1858/2018 (92%)	-0.27	18 (0%) 82 59	70, 112, 159, 208	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	605	ALA	4.2
1	B	606	CYS	3.3
1	B	589	GLU	3.2
1	A	495	TYR	3.1
1	B	557	SER	3.0
1	B	635	GLY	2.7
1	B	686	ASN	2.5
1	A	551	GLY	2.5
1	B	590	LYS	2.5
1	B	598	LYS	2.2
1	B	547	GLY	2.2
1	B	721	ASN	2.2
1	B	639	GLU	2.2
1	B	661	TYR	2.2
1	A	720	ILE	2.1
1	A	877	ASN	2.1
1	A	680	GLU	2.0
1	B	687	ASP	2.0

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

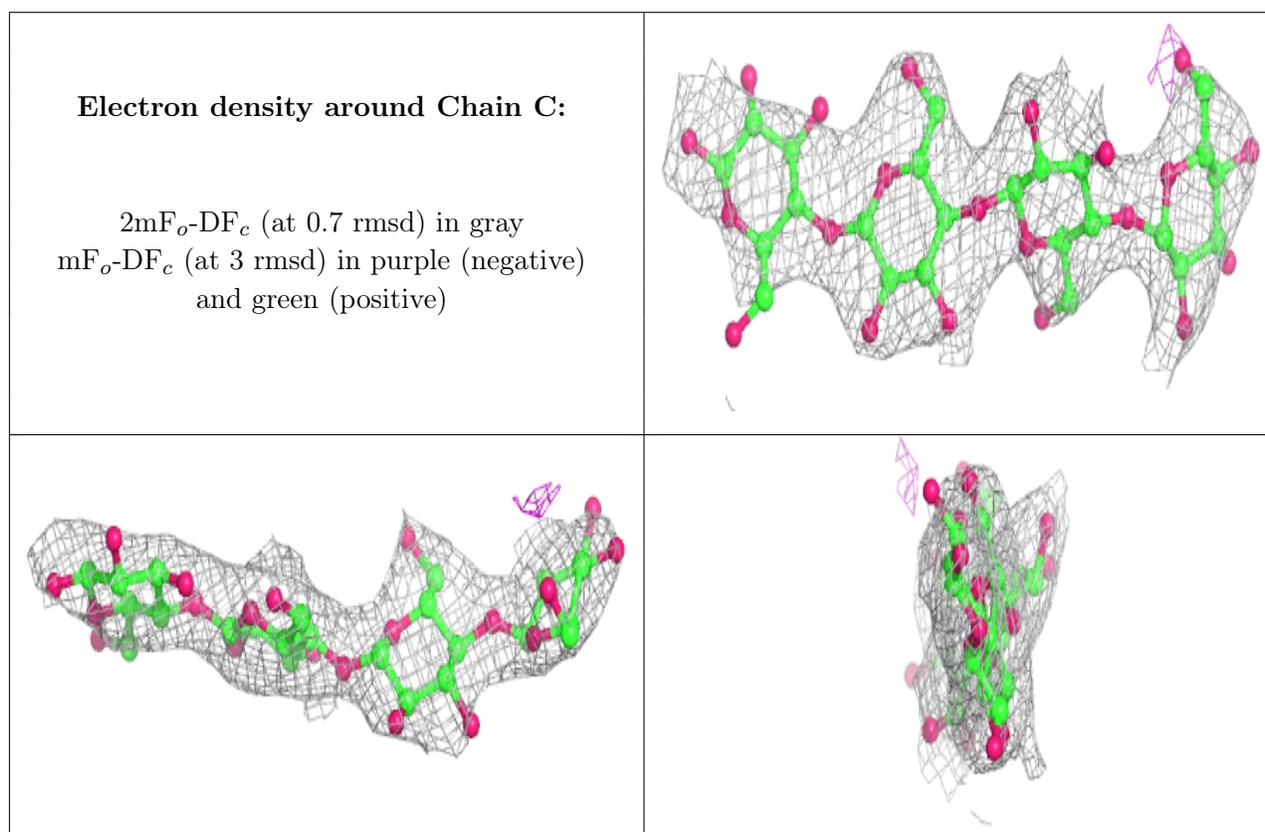
There are no non-standard protein/DNA/RNA residues in this entry.

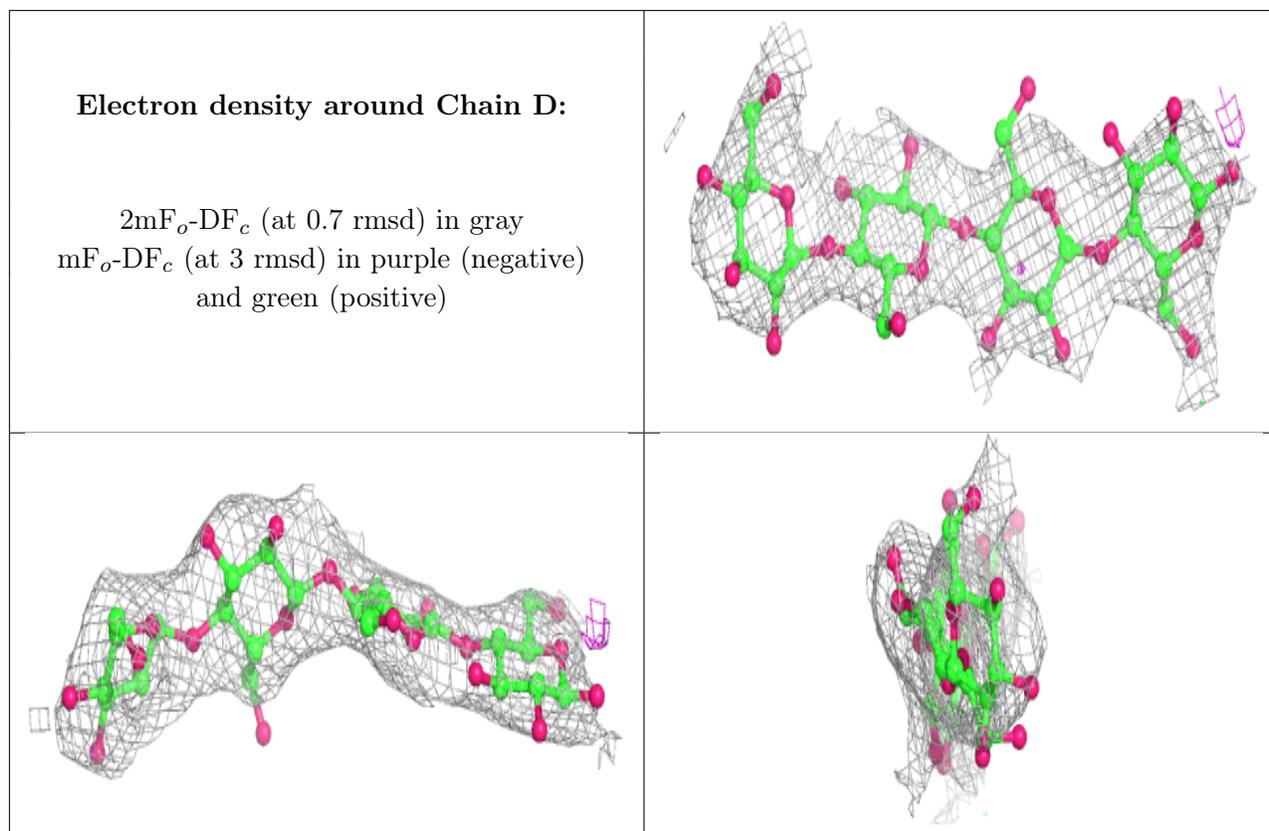
### 6.3 Carbohydrates [i](#)

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	0.79	0.43	143,153,154,156	0
2	BGC	C	3	11/12	0.87	0.20	125,133,140,146	0
2	BGC	C	4	11/12	0.87	0.24	134,139,144,145	0
2	BGC	C	1	12/12	0.87	0.36	128,142,147,153	0
2	BGC	D	4	11/12	0.90	0.20	145,152,156,156	0
2	BGC	D	2	11/12	0.91	0.34	121,130,135,135	0
2	BGC	D	3	11/12	0.94	0.15	124,127,133,139	0
2	BGC	C	2	11/12	0.95	0.17	91,103,112,114	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](#)

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
3	PO4	A	1005	5/5	0.92	0.16	106,111,113,117	0
3	PO4	B	1005	5/5	0.97	0.11	105,106,114,115	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.