wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 13, 2020 – 12:22 AM JST

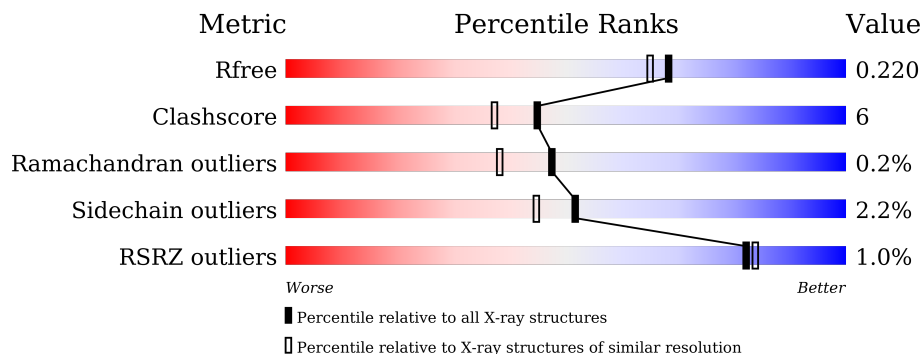|		
PDB ID	:	6LCF
Title	:	Crystal Structure of beta-L-arabinobiose binding protein - native
Authors	:	Miyake, M.; Arakawa, T.; Fushinobu, S.
Deposited on	:	2019-11-18
Resolution	:	1.92 Å(reported)

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1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	Percentile Ranks	Value
Rfree		0.220
Clashscore		6
Ramachandran outliers		0.2%
Sidechain outliers		2.2%
RSRZ outliers		1.0%

Worse ←→ Better

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 >=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 <=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	436	.% 79% 13% · 8%
1	B	436	.% 82% 10% 8%
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 6296 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 ABC transporter substrate binding component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3023	1894	502	620	7			
1	B	403	Total	C	N	O	S	0	0	0
			3023	1894	502	620	7			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	MET	-	expression tag	UNP A0A0A1GL90
A	7	GLY	-	expression tag	UNP A0A0A1GL90
A	8	HIS	-	expression tag	UNP A0A0A1GL90
A	9	HIS	-	expression tag	UNP A0A0A1GL90
A	10	HIS	-	expression tag	UNP A0A0A1GL90
A	11	HIS	-	expression tag	UNP A0A0A1GL90
A	12	HIS	-	expression tag	UNP A0A0A1GL90
A	13	HIS	-	expression tag	UNP A0A0A1GL90
A	14	HIS	-	expression tag	UNP A0A0A1GL90
A	15	HIS	-	expression tag	UNP A0A0A1GL90
A	16	HIS	-	expression tag	UNP A0A0A1GL90
A	17	HIS	-	expression tag	UNP A0A0A1GL90
A	18	SER	-	expression tag	UNP A0A0A1GL90
A	19	SER	-	expression tag	UNP A0A0A1GL90
A	20	GLY	-	expression tag	UNP A0A0A1GL90
A	21	HIS	-	expression tag	UNP A0A0A1GL90
A	22	ILE	-	expression tag	UNP A0A0A1GL90
A	23	GLU	-	expression tag	UNP A0A0A1GL90
A	24	GLY	-	expression tag	UNP A0A0A1GL90
A	25	ARG	-	expression tag	UNP A0A0A1GL90
A	26	HIS	-	expression tag	UNP A0A0A1GL90
A	27	MET	-	expression tag	UNP A0A0A1GL90
B	6	MET	-	expression tag	UNP A0A0A1GL90
B	7	GLY	-	expression tag	UNP A0A0A1GL90
B	8	HIS	-	expression tag	UNP A0A0A1GL90

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Chain	Residue	Modelled	Actual	Comment	Reference
B	9	HIS	-	expression tag	UNP A0A0A1GL90
B	10	HIS	-	expression tag	UNP A0A0A1GL90
B	11	HIS	-	expression tag	UNP A0A0A1GL90
B	12	HIS	-	expression tag	UNP A0A0A1GL90
B	13	HIS	-	expression tag	UNP A0A0A1GL90
B	14	HIS	-	expression tag	UNP A0A0A1GL90
B	15	HIS	-	expression tag	UNP A0A0A1GL90
B	16	HIS	-	expression tag	UNP A0A0A1GL90
B	17	HIS	-	expression tag	UNP A0A0A1GL90
B	18	SER	-	expression tag	UNP A0A0A1GL90
B	19	SER	-	expression tag	UNP A0A0A1GL90
B	20	GLY	-	expression tag	UNP A0A0A1GL90
B	21	HIS	-	expression tag	UNP A0A0A1GL90
B	22	ILE	-	expression tag	UNP A0A0A1GL90
B	23	GLU	-	expression tag	UNP A0A0A1GL90
B	24	GLY	-	expression tag	UNP A0A0A1GL90
B	25	ARG	-	expression tag	UNP A0A0A1GL90
B	26	HIS	-	expression tag	UNP A0A0A1GL90
B	27	MET	-	expression tag	UNP A0A0A1GL90

- Molecule 2 is an oligosaccharide called beta-L-arabinofuranose-(1-2)-beta-L-arabinofuranose.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total 19	C 10	O 9	0	0	0
2	D	2	Total 19	C 10	O 9	0	0	0

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total 109	O 109	0	0
3	B	103	Total 103	O 103	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: ABC transporter substrate binding component
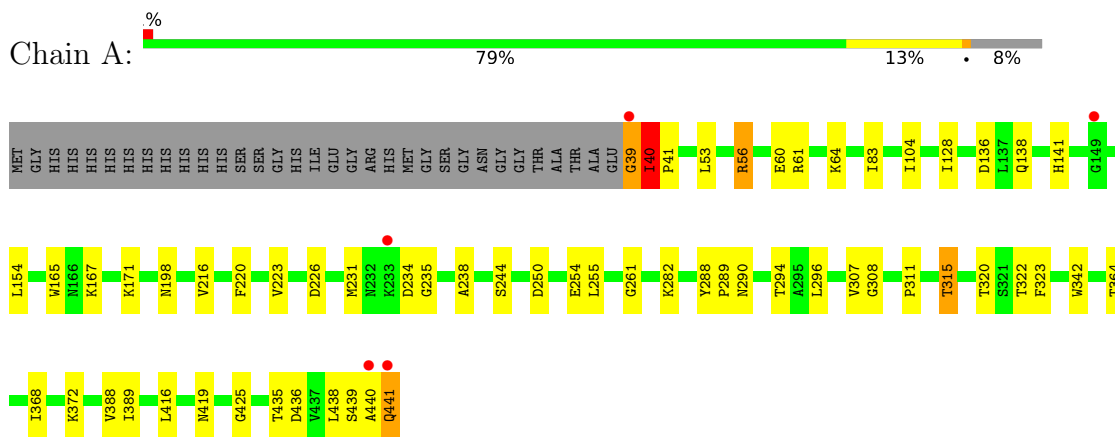
Chain A:

- Molecule 1: ABC transporter substrate binding component
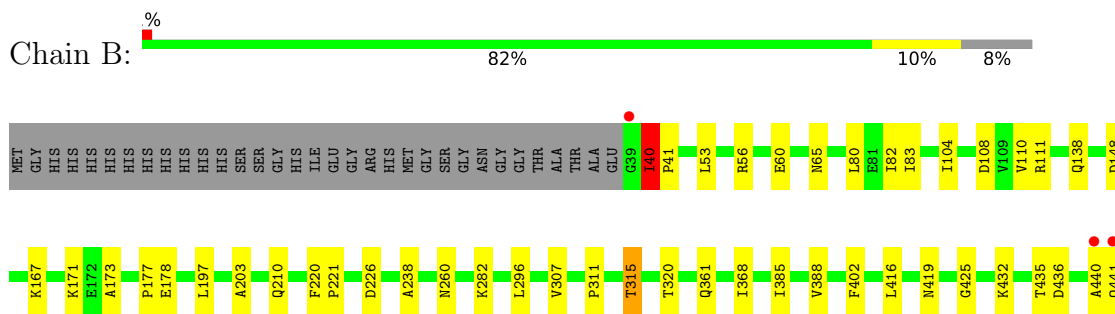
Chain B:

- Molecule 2: beta-L-arabinofuranose-(1-2)-beta-L-arabinofuranose

Chain C:

- Molecule 2: beta-L-arabinofuranose-(1-2)-beta-L-arabinofuranose

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	69.35Å 52.96Å 102.77Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	46.97 – 1.92 46.93 – 1.92	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.97-1.92) 100.0 (46.93-1.92)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.96 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.182 , 0.214 0.191 , 0.220	Depositor DCC
R_{free} test set	2712 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å^2)	15.7	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 39.1	EDS
L-test for twinning[2]	$< |L| > = 0.47, < L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6296	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5   Model quality ⓘ

5.1   Standard geometry ⓘ

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

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	0.73	0/3083	0.84	2/4188 (0.0%)
1	B	0.72	0/3083	0.81	0/4188
All	All	0.72	0/6166	0.83	2/8376 (0.0%)

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

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	40	ILE	CB-CA-C	7.41	126.42	111.60
1	A	56	ARG	NE-CZ-NH1	-5.79	117.41	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	39	GLY	Peptide

5.2   Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3023	0	2919	50	0
1	B	3023	0	2919	25	0
2	C	19	0	0	1	0
2	D	19	0	0	1	0
3	A	109	0	0	1	0
3	B	103	0	0	2	0
All	All	6296	0	5838	77	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.

The worst 5 of 77 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:238:ALA:H	1:B:435:THR:HG22	1.13	1.10
1:A:238:ALA:H	1:A:435:THR:HG22	1.17	1.03
1:A:216:VAL:HG23	1:A:220:PHE:CD2	2.01	0.96
1:A:231:MET:HE3	1:A:438:LEU:HD13	1.45	0.95
1:A:216:VAL:HG23	1:A:220:PHE:CE2	2.01	0.95

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	401/436 (92%)	390 (97%)	10 (2%)	1 (0%)	47	38
1	B	401/436 (92%)	391 (98%)	9 (2%)	1 (0%)	47	38
All	All	802/872 (92%)	781 (97%)	19 (2%)	2 (0%)	47	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	ILE
1	B	40	ILE

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	318/342 (93%)	313 (98%)	5 (2%)	62	58
1	B	318/342 (93%)	309 (97%)	9 (3%)	43	34
All	All	636/684 (93%)	622 (98%)	14 (2%)	52	45

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	60	GLU
1	B	138	GLN
1	B	282	LYS
1	B	40	ILE
1	B	210	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	FUB	C	1	2	10,10,10	1.39	2 (20%)	13,14,14	2.56	7 (53%)
2	FUB	C	2	2	9,9,10	0.57	0	10,12,14	1.14	1 (10%)
2	FUB	D	1	2	10,10,10	3.47	3 (30%)	13,14,14	2.78	7 (53%)
2	FUB	D	2	2	9,9,10	2.25	2 (22%)	10,12,14	1.99	2 (20%)

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	FUB	C	1	2	-	2/2/18/18	0/1/1/1
2	FUB	C	2	2	-	0/2/15/18	0/1/1/1
2	FUB	D	1	2	-	2/2/18/18	0/1/1/1
2	FUB	D	2	2	-	0/2/15/18	0/1/1/1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	FUB	C1-C2	7.11	1.61	1.52
2	D	1	FUB	O2-C2	-6.96	1.26	1.43
2	D	2	FUB	O2-C2	6.02	1.56	1.43
2	D	1	FUB	O4-C4	3.82	1.53	1.45
2	C	1	FUB	O4-C4	2.86	1.51	1.45

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	FUB	O5-C5-C4	-4.97	94.25	111.29
2	D	1	FUB	O2-C2-C3	4.40	126.04	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	1	FUB	O5-C5-C4	-4.13	97.11	111.29
2	D	1	FUB	O1-C1-C2	4.05	129.71	110.39
2	D	2	FUB	O2-C2-C1	-3.99	99.07	110.97

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	FUB	O4-C4-C5-O5
2	D	1	FUB	O4-C4-C5-O5
2	C	1	FUB	C3-C4-C5-O5
2	D	1	FUB	C3-C4-C5-O5

There are no ring outliers.

2 monomers are involved in 2 short contacts:

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

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain D	
Bond lengths	Bond angles
Torsions	Rings
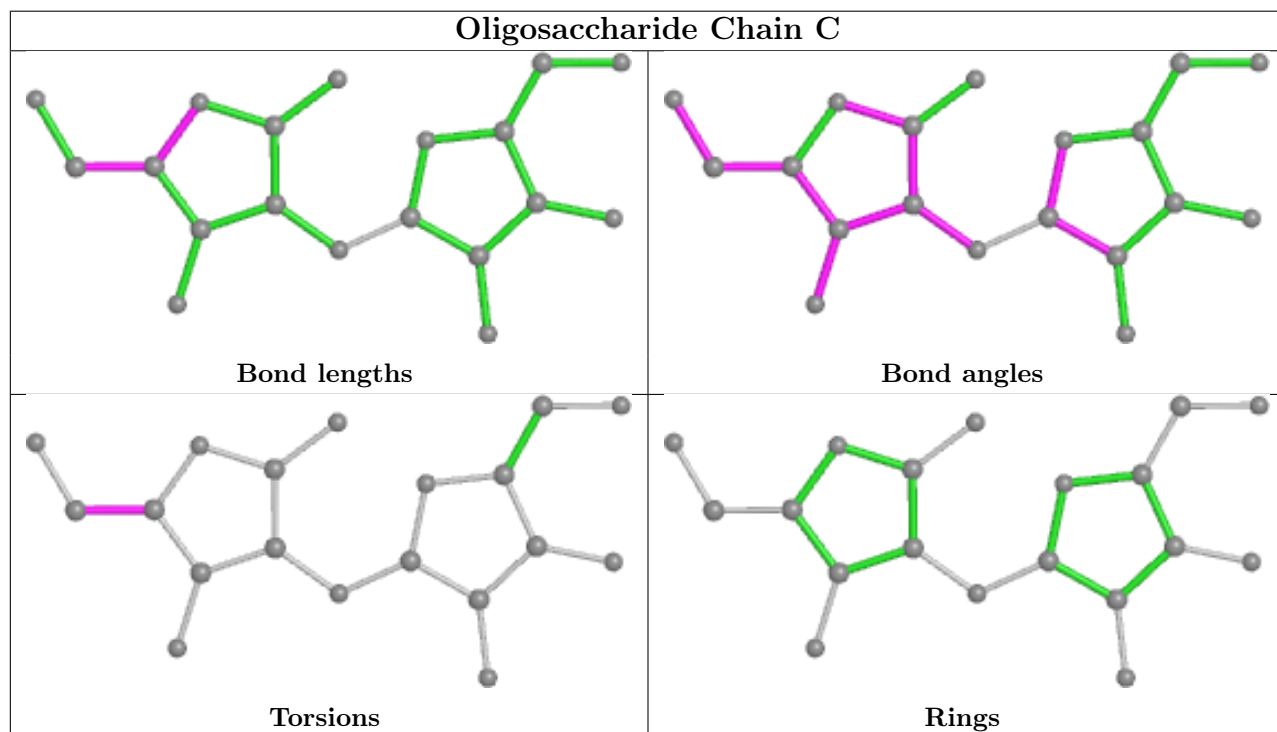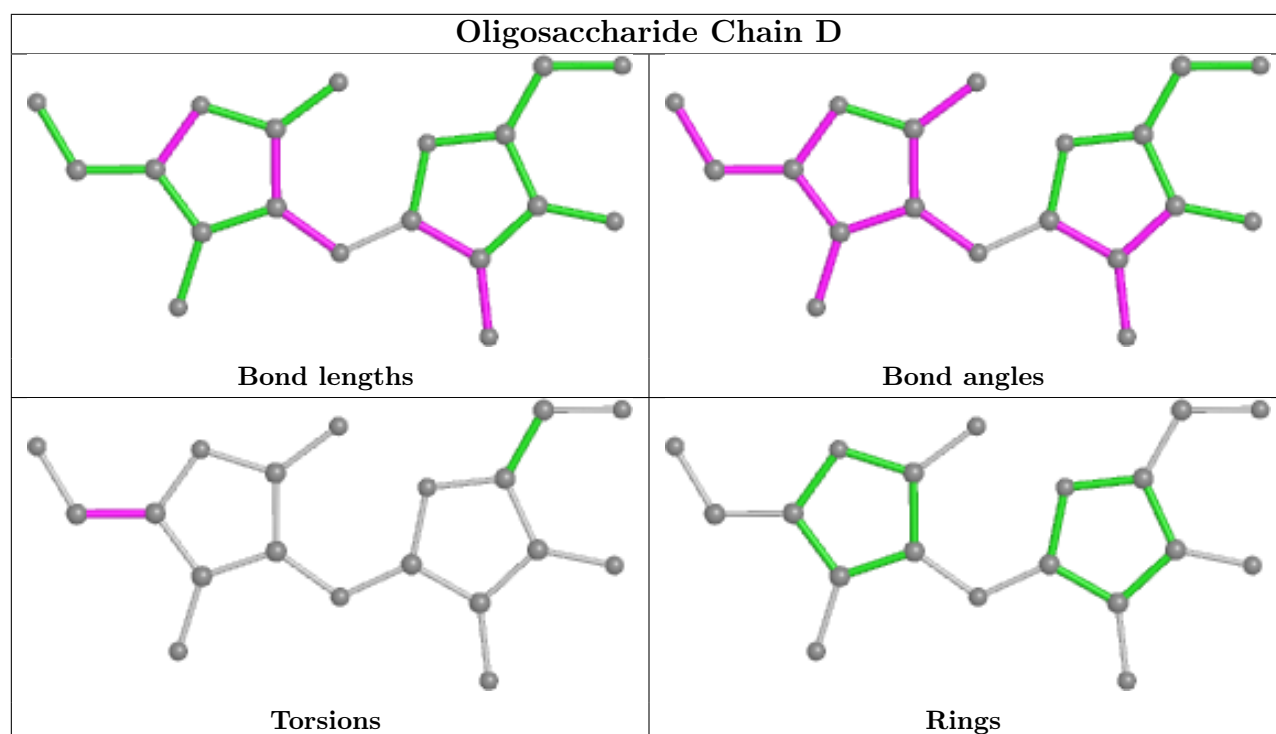
5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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^2)	Q<0.9
1	A	403/436 (92%)	-0.27	5 (1%)	79	81	11, 18, 31, 48	0
1	B	403/436 (92%)	-0.28	3 (0%)	87	89	11, 18, 34, 65	0
All	All	806/872 (92%)	-0.27	8 (0%)	82	84	11, 18, 31, 65	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	39	GLY	6.2
1	A	440	ALA	3.8
1	B	440	ALA	3.7
1	B	39	GLY	3.4
1	B	441	GLN	2.8

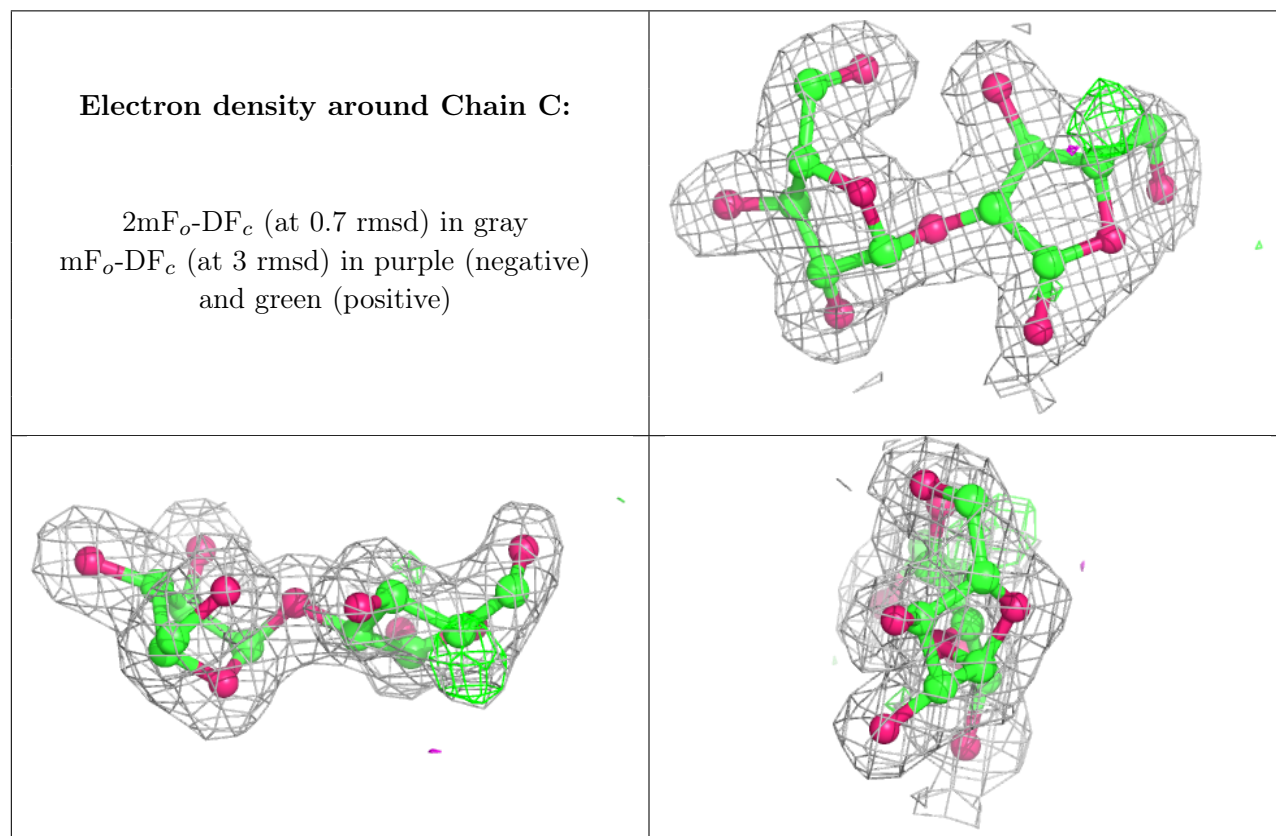## 6.2   Non-standard residues in protein, DNA, RNA chains ⓘ

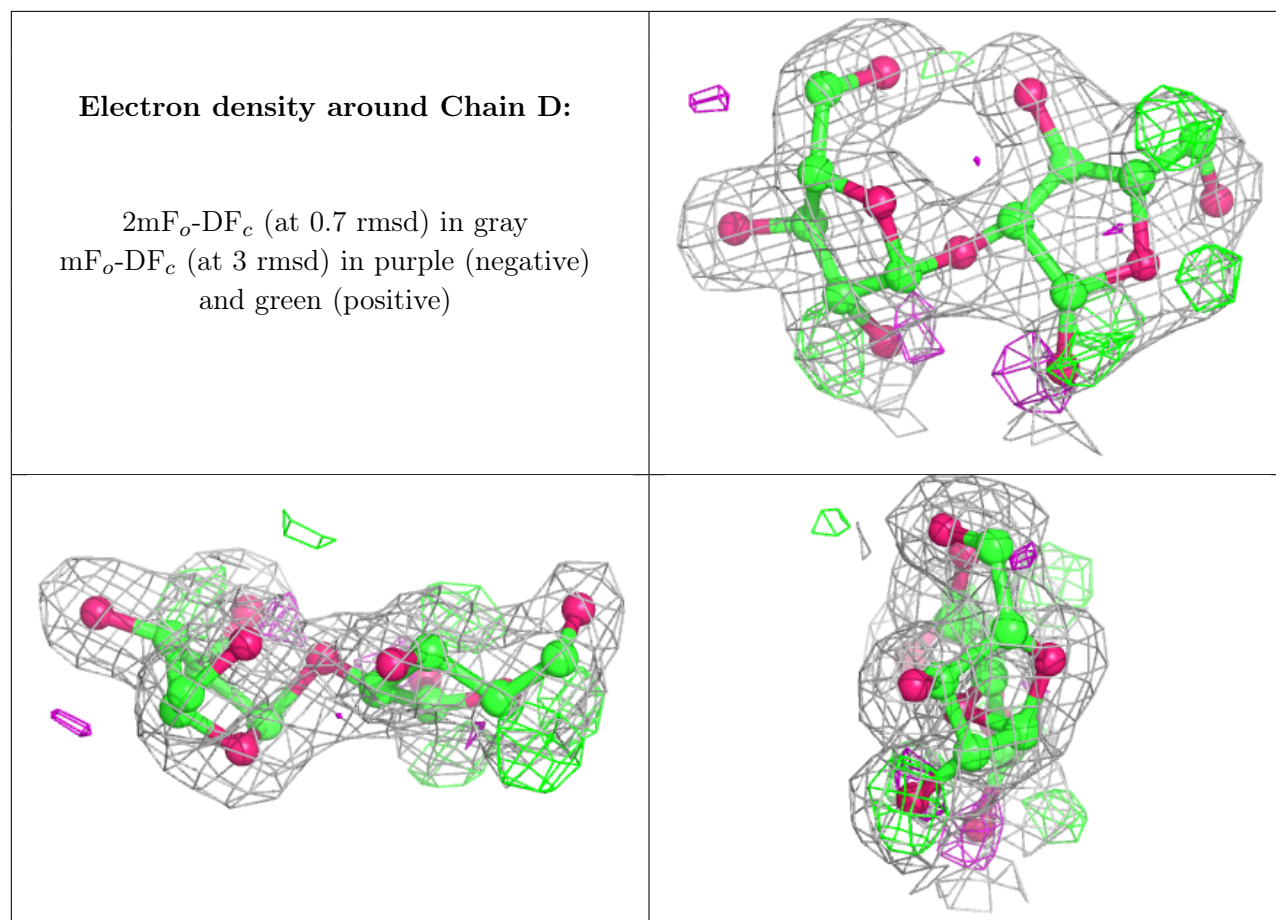There are no non-standard protein/DNA/RNA residues in this entry.

6.3   Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th 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^2)	Q<0.9
2	FUB	D	1	10/10	0.85	0.17	14,17,23,25	0
2	FUB	D	2	9/10	0.91	0.10	11,12,13,13	0
2	FUB	C	1	10/10	0.92	0.11	14,17,20,22	0
2	FUB	C	2	9/10	0.98	0.07	11,12,12,13	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.

Electron density around Chain C:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

Electron density around Chain D:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative) and green (positive)

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.