



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 08:20 AM BST

PDB ID : 4XUR
Title : Structure of the CBM22-2 xylan-binding domain from *Paenibacillus barcinonensis* Xyn10C in complex with xylotetraose
Authors : Sainz-Polo, M.A.; Sanz-Aparicio, J.
Deposited on : 2015-01-26
Resolution : 1.67 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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 : 2.13.1
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) : 2.13.1

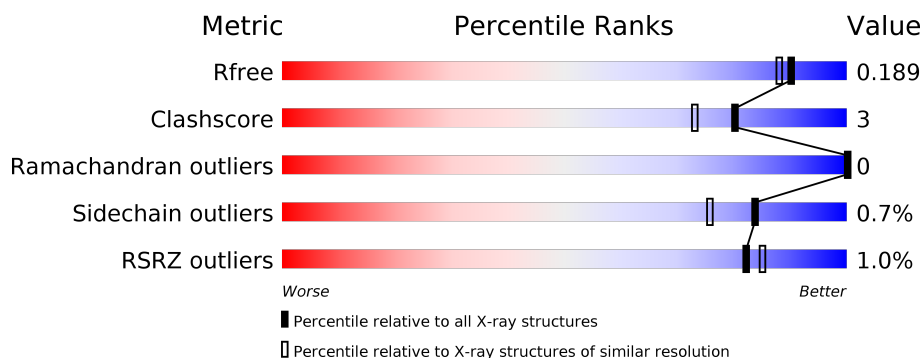
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	6780 (1.70-1.66)
Clashscore	141614	7310 (1.70-1.66)
Ramachandran outliers	138981	7173 (1.70-1.66)
Sidechain outliers	138945	7172 (1.70-1.66)
RSRZ outliers	127900	6661 (1.70-1.66)

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 on 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 $\leq 5\%$. 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	181	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>7%</div> <div>••</div> <div>11%</div> </div> </div>
1	B	181	<div> <div></div> <div> <div>86%</div> <div>•</div> <div>11%</div> </div> </div>
1	C	181	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>5%</div> <div>11%</div> </div> </div>
2	D	4	<div> <div></div> <div>100%</div> </div>
3	E	2	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>
4	F	3	<div> <div></div> <div> <div>33%</div> <div>67%</div> </div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3987 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 Endo-1,4-beta-xylanase C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1263	796	211	253	3			
1	B	161	Total	C	N	O	S	0	0	0
			1263	796	211	253	3			
1	C	161	Total	C	N	O	S	0	0	0
			1263	796	211	253	3			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	177	ASP	TYR	conflict	UNP O69230
B	177	ASP	TYR	conflict	UNP O69230
C	177	ASP	TYR	conflict	UNP O69230

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



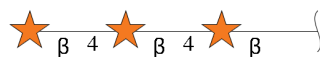
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	4	Total	C	O	0	0	0
			37	20	17			

- Molecule 3 is an oligosaccharide called beta-D-xylopyranose-(1-4)-beta-D-xylopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	E	2	Total	C	O	0	0	0
			18	10	8			

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

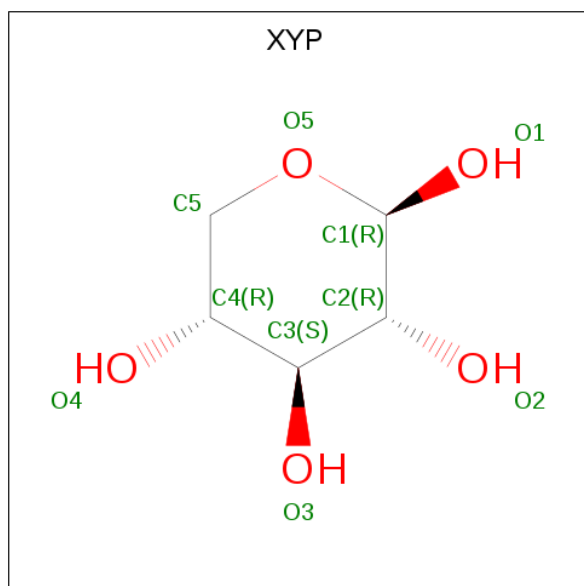


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	F	3	Total	C	O	0	0	0
			28	15	13			

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 6 is beta-D-xylopyranose (three-letter code: XYP) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			9	5	4		
6	C	1	Total	C	O	0	0
			9	5	4		

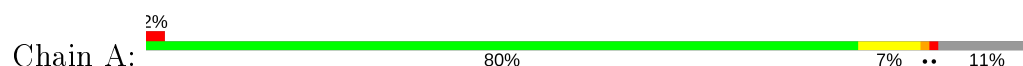
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	33	Total	O	0	0
			33	33		
7	B	30	Total	O	0	0
			30	30		
7	C	31	Total	O	0	0
			31	31		

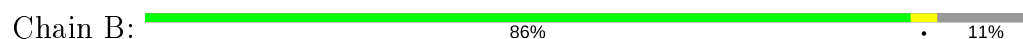
3 Residue-property plots [i](#)

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.

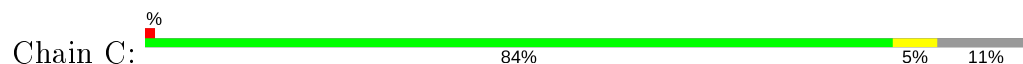
- Molecule 1: Endo-1,4-beta-xylanase C



- Molecule 1: Endo-1,4-beta-xylanase C



- Molecule 1: Endo-1,4-beta-xylanase C



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



- Molecule 3: beta-D-xylopyranose-(1-4)-beta-D-xylopyranose



- Molecule 4: beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain F:

33%

67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	92.77Å 92.77Å 48.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.55 – 1.67 33.55 – 1.67	Depositor EDS
% Data completeness (in resolution range)	100.0 (33.55-1.67) 100.0 (33.55-1.67)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 1.67Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.158 , 0.194 0.163 , 0.189	Depositor DCC
R_{free} test set	2769 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 17.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.124 for -h,-k,l 0.119 for h,-h-k,-l 0.456 for -k,-h,-l	Xtriage
Reported twinning fraction	0.537 for H, K, L 0.463 for -K, -H, -L	Depositor
Outliers	0 of 54477 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3987	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.19% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.02	2/1291 (0.2%)	0.94	3/1753 (0.2%)
1	B	0.94	0/1291	0.86	0/1753
1	C	0.97	1/1291 (0.1%)	0.83	1/1753 (0.1%)
All	All	0.98	3/3873 (0.1%)	0.88	4/5259 (0.1%)

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	267	GLU	CD-OE2	-6.74	1.18	1.25
1	A	308	TRP	CB-CG	-6.16	1.39	1.50
1	C	267	GLU	CD-OE2	-5.65	1.19	1.25

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	301	ARG	NE-CZ-NH1	6.41	123.50	120.30
1	C	287	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	187	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	A	309	LYS	C-N-CA	5.01	134.23	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	308	TRP	Peptide

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	1263	0	1216	16	0
1	B	1263	0	1216	2	0
1	C	1263	0	1216	7	0
2	D	37	0	0	0	0
3	E	18	0	0	0	0
4	F	28	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
6	B	9	0	0	0	0
6	C	9	0	0	0	0
7	A	33	0	0	0	0
7	B	30	0	0	0	0
7	C	31	0	0	0	0
All	All	3987	0	3648	20	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 3.

All (20) 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:316:GLU:OE1	1:C:272:ALA:HB1	1.90	0.71
1:A:216:THR:HG21	1:A:309:LYS:HA	1.76	0.68
1:A:316:GLU:HB2	1:C:274:THR:HG21	1.79	0.64
1:A:309:LYS:HB3	1:A:310:GLN:HA	1.82	0.61
1:A:310:GLN:O	1:A:313:GLU:N	2.32	0.59
1:A:309:LYS:HG2	1:A:314:VAL:CG2	2.33	0.57
1:A:193:LYS:NZ	1:A:195:GLU:OE1	2.37	0.57
1:A:309:LYS:HG2	1:A:314:VAL:HG22	1.89	0.55
1:A:316:GLU:CB	1:C:274:THR:HG21	2.38	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ASN:HB2	1:A:280:GLN:HG2	1.91	0.52
1:A:308:TRP:CE3	1:A:309:LYS:HB2	2.46	0.50
1:B:252:ASN:OD1	1:B:273:SER:HB2	2.12	0.50
1:A:216:THR:HG23	1:A:309:LYS:HG3	1.95	0.49
1:C:225:LEU:N	1:C:226:PRO:CD	2.75	0.49
1:A:316:GLU:CD	1:C:272:ALA:HB1	2.34	0.48
1:A:308:TRP:CZ3	1:A:309:LYS:HE2	2.50	0.47
1:C:236:ILE:HG12	1:C:329:ILE:HG12	1.98	0.46
1:A:242:TYR:CD2	1:A:320:ILE:HD13	2.53	0.43
1:B:236:ILE:HD11	1:B:288:ILE:CD1	2.49	0.42
1:A:316:GLU:OE2	1:C:272:ALA:HB1	2.20	0.41

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	159/181 (88%)	156 (98%)	3 (2%)	0	100	100
1	B	159/181 (88%)	157 (99%)	2 (1%)	0	100	100
1	C	159/181 (88%)	157 (99%)	2 (1%)	0	100	100
All	All	477/543 (88%)	470 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	137/150 (91%)	136 (99%)	1 (1%)	84	76
1	B	137/150 (91%)	136 (99%)	1 (1%)	84	76
1	C	137/150 (91%)	136 (99%)	1 (1%)	84	76
All	All	411/450 (91%)	408 (99%)	3 (1%)	84	76

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	301	ARG
1	B	301	ARG
1	C	301	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	HIS
1	A	277	ASN
1	A	280	GLN
1	A	310	GLN
1	B	185	ASN
1	B	203	ASN
1	B	277	ASN
1	B	280	GLN
1	C	229	GLN
1	C	260	ASN
1	C	277	ASN
1	C	280	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 ⓘ

9 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	XYP	D	1	2	10,10,10	1.04	0	14,14,14	1.90	2 (14%)
2	XYP	D	2	2	9,9,10	1.30	1 (11%)	10,12,14	1.01	1 (10%)
2	XYP	D	3	2	9,9,10	1.10	1 (11%)	10,12,14	1.13	0
2	XYP	D	4	2	9,9,10	1.30	2 (22%)	10,12,14	0.67	0
3	XYP	E	1	3	9,9,10	0.96	0	10,12,14	1.18	2 (20%)
3	XYP	E	2	3	9,9,10	0.47	0	10,12,14	0.91	0
4	XYP	F	1	4	10,10,10	0.93	0	14,14,14	1.28	2 (14%)
4	XYP	F	2	4	9,9,10	1.16	2 (22%)	10,12,14	1.42	2 (20%)
4	XYP	F	3	4	9,9,10	0.82	0	10,12,14	0.70	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	XYP	D	1	2	-	-	0/1/1/1
2	XYP	D	2	2	-	-	0/1/1/1
2	XYP	D	3	2	-	-	0/1/1/1
2	XYP	D	4	2	-	-	0/1/1/1
3	XYP	E	1	3	-	-	0/1/1/1
3	XYP	E	2	3	-	-	0/1/1/1
4	XYP	F	1	4	-	-	0/1/1/1
4	XYP	F	2	4	-	-	0/1/1/1
4	XYP	F	3	4	-	-	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	2	XYP	C4-C3	2.31	1.55	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3	XYP	O5-C1	2.27	1.47	1.42
2	D	4	XYP	O5-C1	2.18	1.47	1.42
2	D	4	XYP	C2-C3	-2.07	1.49	1.52
2	D	2	XYP	O2-C2	2.05	1.47	1.43
4	F	2	XYP	O3-C3	-2.04	1.38	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	XYP	C5-C4-C3	5.38	116.28	109.67
4	F	1	XYP	C5-C4-C3	2.94	113.28	109.67
4	F	2	XYP	C5-C4-C3	2.64	112.92	109.67
2	D	1	XYP	C1-C2-C3	2.59	115.68	110.31
2	D	2	XYP	C1-C2-C3	2.24	112.41	109.67
3	E	1	XYP	C1-C2-C3	2.22	112.40	109.67
3	E	1	XYP	C4-C3-C2	-2.15	108.37	110.92
4	F	1	XYP	O5-C1-C2	2.14	112.61	109.43
4	F	2	XYP	C5-O5-C1	2.02	114.62	111.52

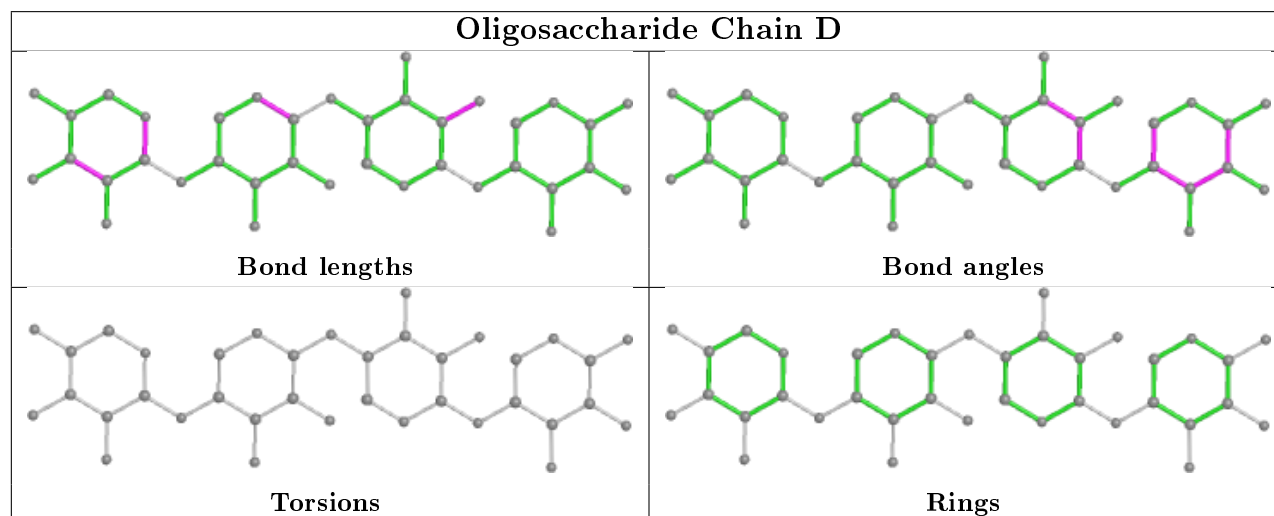
There are no chirality outliers.

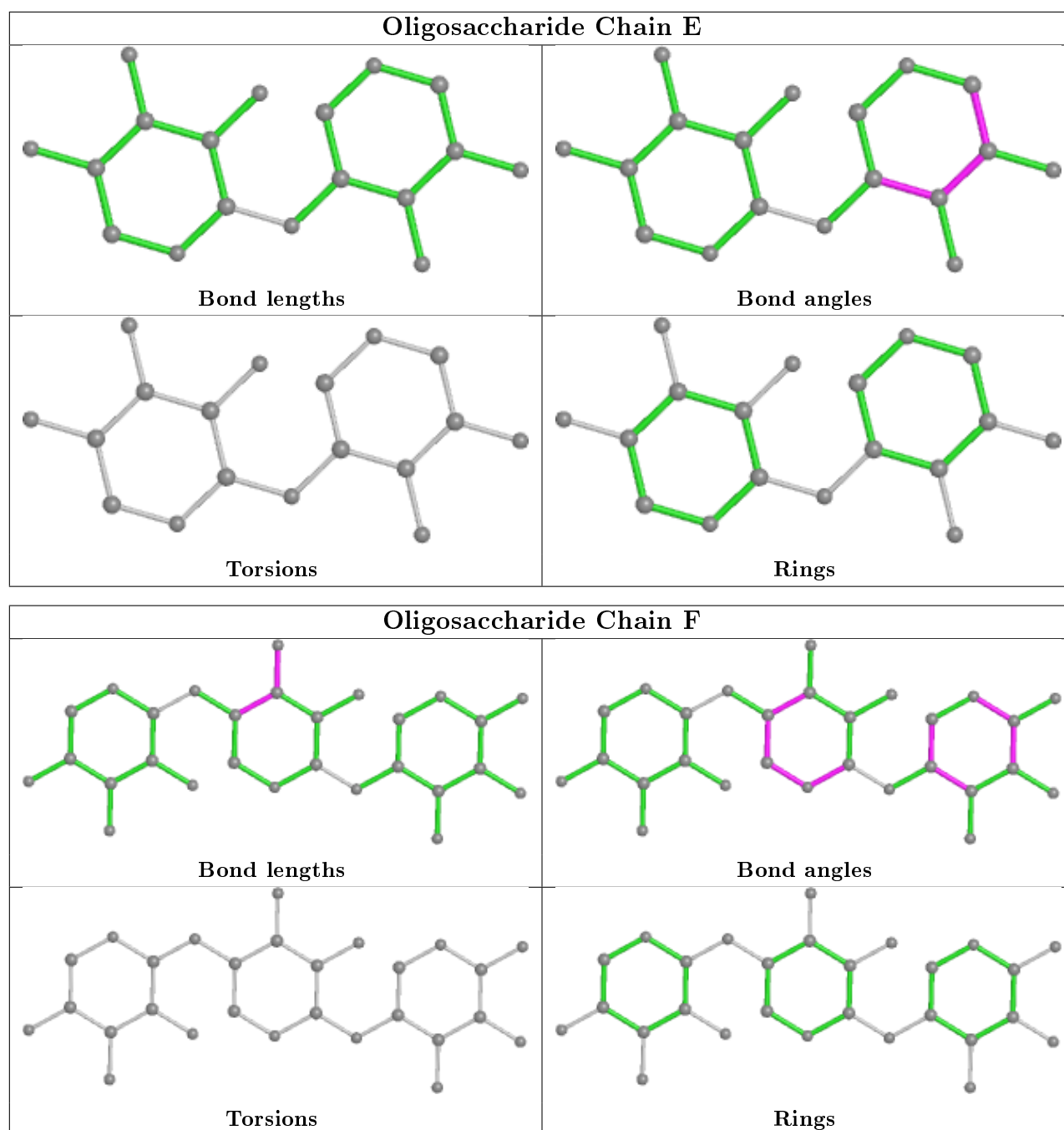
There are no torsion outliers.

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 ⓘ

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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
6	XYP	C	1004	-	9,9,10	0.61	0	10,12,14	1.32	2 (20%)
6	XYP	B	1003	-	9,9,10	0.95	0	10,12,14	1.58	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
6	XYP	C	1004	-	-	-	0/1/1/1
6	XYP	B	1003	-	-	-	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(°)
6	B	1003	XYP	C1-C2-C3	-3.19	105.74	109.67
6	B	1003	XYP	C5-C4-C3	2.63	112.90	109.67
6	C	1004	XYP	C1-C2-C3	-2.48	106.62	109.67
6	C	1004	XYP	C5-C4-C3	2.37	112.58	109.67

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

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 [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, 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	161/181 (88%)	-0.30	3 (1%) 66 70	6, 14, 28, 57	0
1	B	161/181 (88%)	-0.38	0 100 100	6, 14, 27, 34	0
1	C	161/181 (88%)	-0.22	2 (1%) 79 82	9, 15, 30, 38	0
All	All	483/543 (88%)	-0.30	5 (1%) 82 85	6, 15, 30, 57	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	SER	3.4
1	A	310	GLN	3.1
1	C	172	ALA	2.9
1	A	314	VAL	2.1
1	C	312	SER	2.1

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, 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.

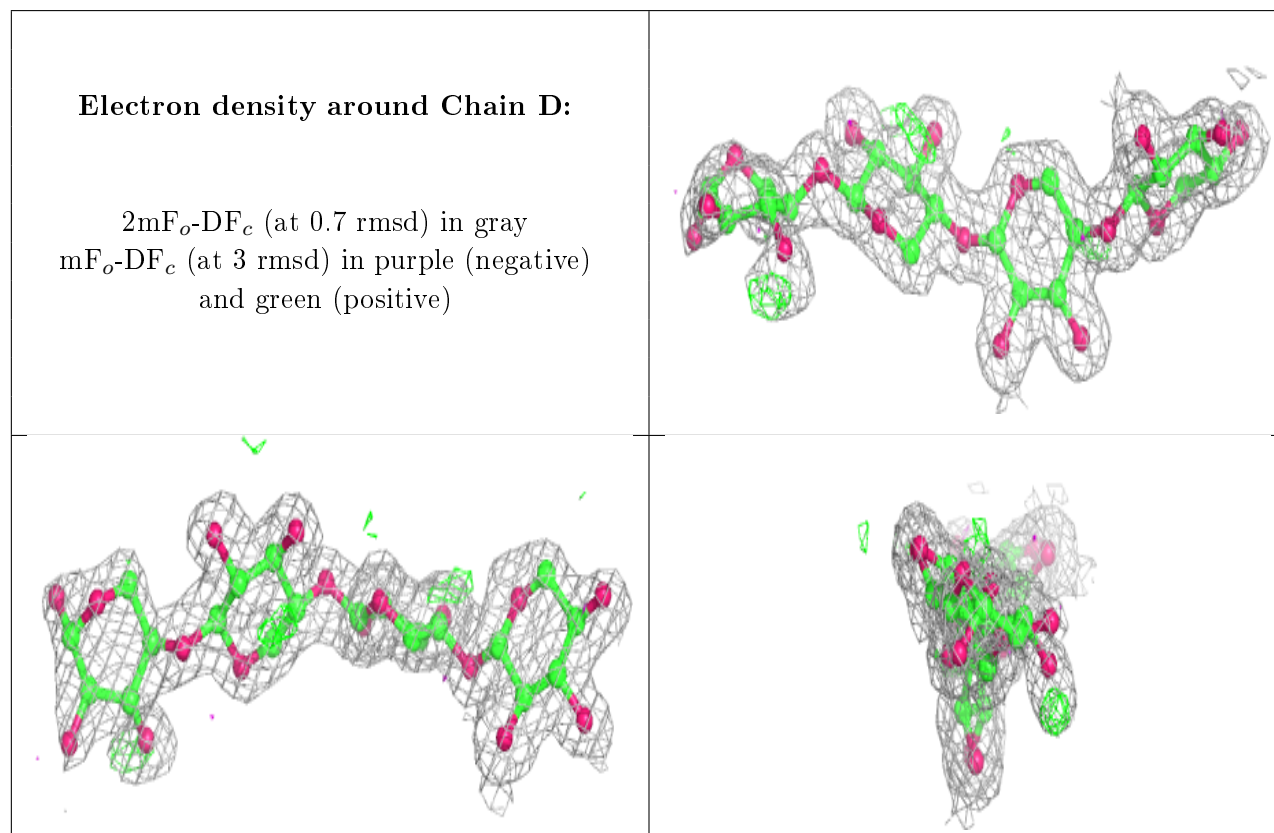
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	XYP	D	1	10/10	0.70	0.20	35,43,47,49	0
4	XYP	F	1	10/10	0.79	0.22	33,37,41,42	0
3	XYP	E	1	9/10	0.85	0.12	22,24,27,28	0

Continued on next page...

Continued from previous page...

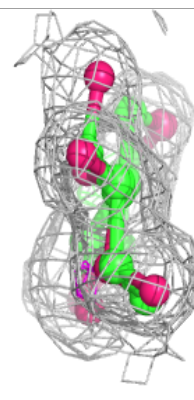
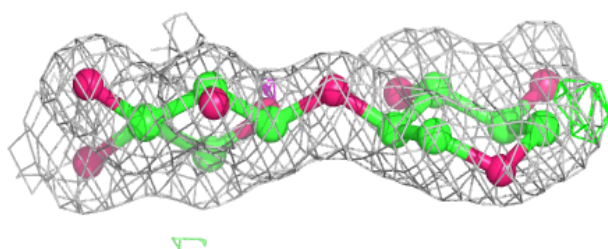
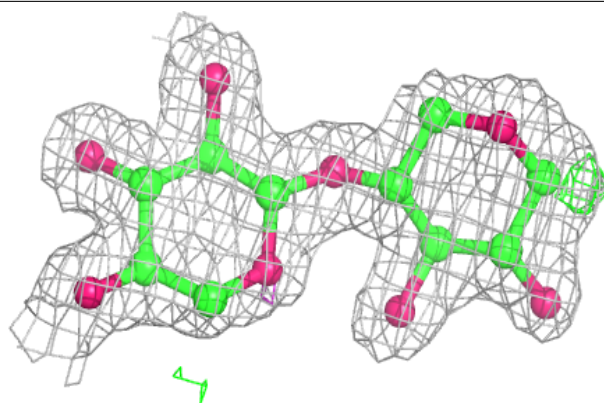
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	XYP	D	2	9/10	0.89	0.10	16,18,20,20	0
4	XYP	F	3	9/10	0.89	0.11	15,16,18,20	0
4	XYP	F	2	9/10	0.89	0.09	19,21,23,23	0
2	XYP	D	4	9/10	0.89	0.11	17,22,24,25	0
3	XYP	E	2	9/10	0.92	0.09	19,20,22,23	0
2	XYP	D	3	9/10	0.93	0.09	13,15,16,17	0

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.

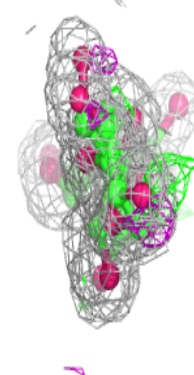
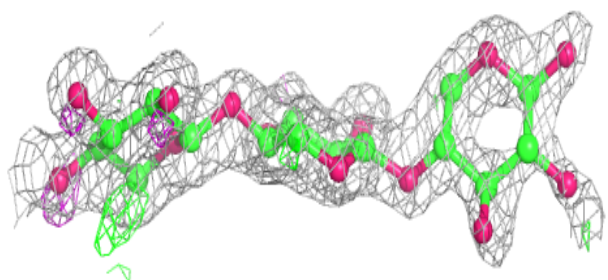
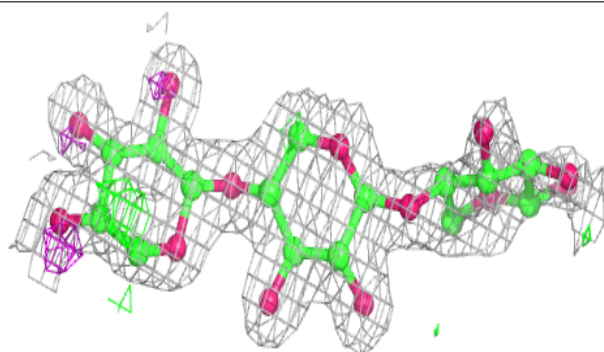


Electron density around Chain E:

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

**Electron density around Chain F:**

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



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, 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(\AA^2)	Q<0.9
6	XYP	C	1004	9/10	0.86	0.11	26,28,31,32	0
6	XYP	B	1003	9/10	0.91	0.11	25,30,33,33	0
5	CA	C	1005	1/1	0.99	0.03	24,24,24,24	0
5	CA	B	1004	1/1	0.99	0.05	14,14,14,14	0
5	CA	A	1005	1/1	0.99	0.03	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.