



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:38 PM UTC

PDB ID : 1UMZ / pdb_00001umz
Title : Xyloglucan endotransglycosylase in complex with the xyloglucan nonasaccharide XLLG.
Authors : Johansson, P.; Brumer, H.; Kallas, A.M.; Henriksson, H.; Denman, S.E.; Teeri, T.T.; Jones, T.A.
Deposited on : 2003-09-03
Resolution : 1.80 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

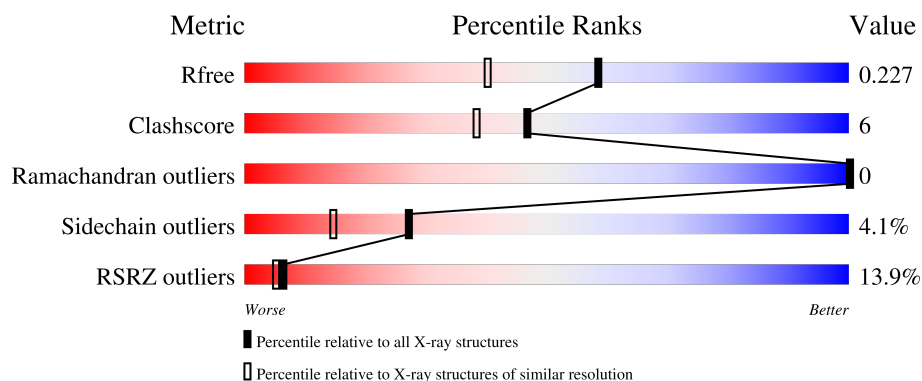
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 $\leq 5\%$. 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	278	18% 81%14% . .
1	B	278	9% 82%10% . 5%
2	C	3	33%33%33%
2	E	3	33%33%33%
3	D	6	33%50%17%

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Mol	Chain	Length	Quality of chain
3	F	6	17% 67% 17%

2 Entry composition [i](#)

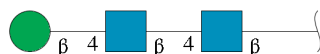
There are 4 unique types of molecules in this entry. The entry contains 4873 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 XYLOGLUCAN ENDOTRANSGLYCOSYLASE.

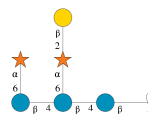
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	15	0	0
			2220	1432	373	404	11			
1	B	264	Total	C	N	O	S	8	1	0
			2206	1426	371	398	11			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	E	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is an oligosaccharide called alpha-D-xylopyranose-(1-6)-beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	6	Total	C	O	0	0	0
			62	34	28			
3	F	6	Total	C	O	0	0	0
			62	34	28			

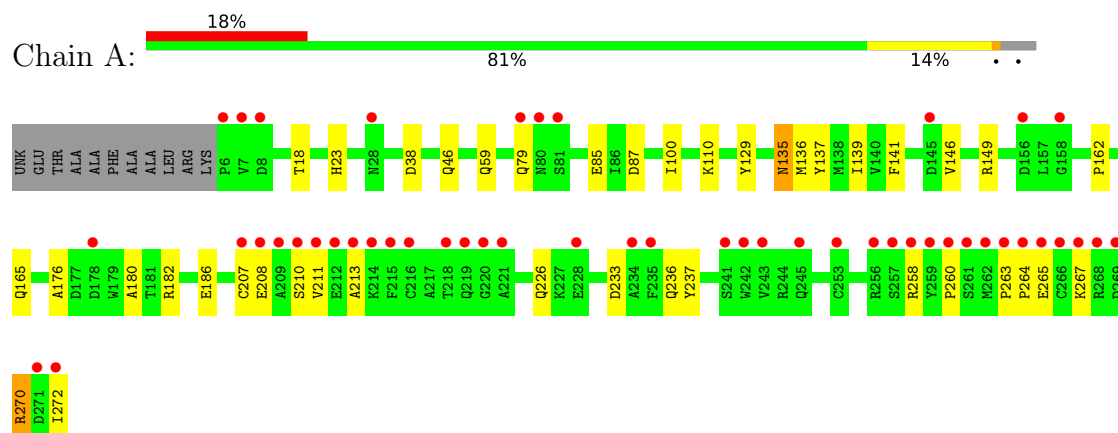
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total 109	O 109	0	0
4	B	136	Total 136	O 136	0	0

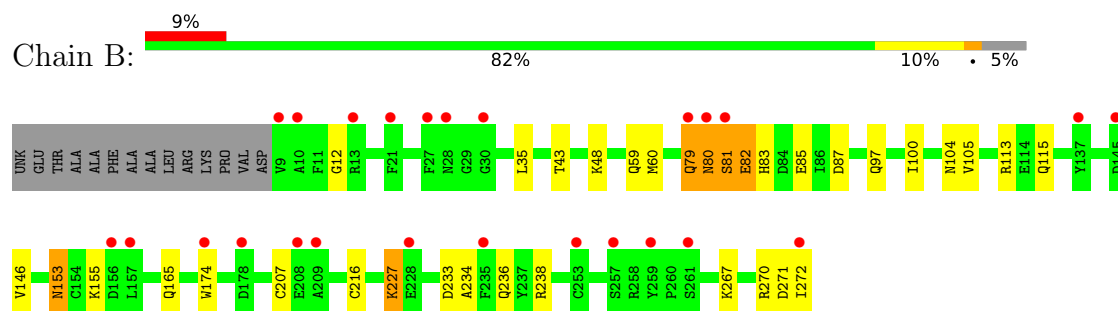
3 Residue-property plots [i](#)

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.

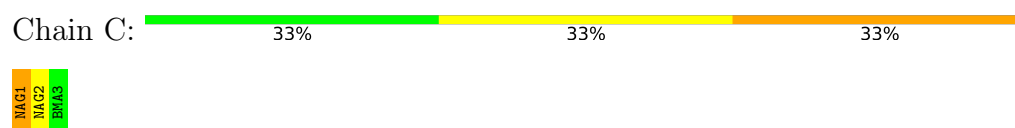
- Molecule 1: XYLOGLUCAN ENDOTRANSGLYCOSYLASE



- Molecule 1: XYLOGLUCAN ENDOTRANSGLYCOSYLASE



- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

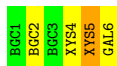


- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 3: alpha-D-xylopyranose-(1-6)-beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



- Molecule 3: alpha-D-xylopyranose-(1-6)-beta-D-glucopyranose-(1-4)-[beta-D-galactopyranose-(1-2)-alpha-D-xylopyranose-(1-6)]beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	188.16 Å 188.16 Å 45.90 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.99 – 1.80 28.99 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.7 (28.99-1.80) 96.7 (28.99-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.80 Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.201 , 0.222 0.208 , 0.227	Depositor DCC
R_{free} test set	4216 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4873	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% 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: NAG, GAL, BMA, BGC, XYS

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	0.79	1/2293 (0.0%)	0.99	9/3104 (0.3%)
1	B	0.77	1/2285 (0.0%)	0.93	5/3095 (0.2%)
All	All	0.78	2/4578 (0.0%)	0.96	14/6199 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	60	MET	SD-CE	-5.04	1.67	1.79
1	A	136	MET	SD-CE	-5.00	1.67	1.79

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	GLN	OE1-CD-NE2	-10.25	112.35	122.60
1	A	226	GLN	OE1-CD-NE2	-9.60	113.00	122.60
1	B	165	GLN	OE1-CD-NE2	-9.41	113.19	122.60
1	B	79	GLN	OE1-CD-NE2	-8.66	113.94	122.60
1	A	211	VAL	N-CA-C	-7.90	102.74	110.72
1	A	165	GLN	CG-CD-NE2	7.08	127.02	116.40
1	A	226	GLN	CG-CD-NE2	6.96	126.84	116.40
1	A	180	ALA	N-CA-C	6.57	118.09	111.07
1	B	165	GLN	CG-CD-NE2	6.54	126.21	116.40
1	B	79	GLN	CG-CD-NE2	6.50	126.15	116.40
1	A	146	VAL	N-CA-C	6.31	114.13	107.76
1	A	176	ALA	CA-C-N	5.80	128.05	120.28
1	A	176	ALA	C-N-CA	5.80	128.05	120.28
1	B	146	VAL	N-CA-C	5.41	113.22	107.76

There are no chirality outliers.

There are no planarity outliers.

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	2220	0	2083	19	0
1	B	2206	0	2070	25	0
2	C	39	0	34	1	0
2	E	39	0	34	3	0
3	D	62	0	52	2	0
3	F	62	0	51	2	0
4	A	109	0	0	1	0
4	B	136	0	0	4	0
All	All	4873	0	4324	48	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.

All (48) 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:80:ASN:OD1	1:B:82:GLU:N	1.88	1.06
1:B:97:GLN:NE2	2:E:1:NAG:O6	1.92	1.03
1:A:260:PRO:O	4:A:301:HOH:O	1.98	0.81
1:B:80:ASN:OD1	1:B:81:SER:N	2.17	0.77
1:B:233:ASP:H	1:B:236:GLN:HE21	1.32	0.76
3:F:2:BGC:C6	3:F:5:XYS:C1	2.65	0.75
1:A:139:ILE:HD11	1:A:162:PRO:CG	2.17	0.74
1:B:12:GLY:O	1:B:48:LYS:NZ	2.21	0.73
1:B:207:CYS:HG	1:B:216:CYS:HG	0.74	0.73
3:D:2:BGC:H6	3:D:5:XYS:C1	1.99	0.72
1:A:139:ILE:HD11	1:A:162:PRO:HG2	1.73	0.70
1:B:174[B]:TRP:CD1	4:B:312:HOH:O	2.45	0.70
1:B:80:ASN:O	1:B:83:HIS:HB3	1.95	0.66
1:A:135:ASN:HD22	1:A:135:ASN:C	2.09	0.61
1:A:23:HIS:HE1	1:A:38:ASP:OD2	1.84	0.60
1:B:97:GLN:NE2	2:E:1:NAG:C6	2.65	0.60
1:B:104:ASN:HD21	1:B:113:ARG:H	1.50	0.59
1:B:153:ASN:HD22	1:B:153:ASN:C	2.10	0.58
1:A:18:THR:HG21	1:A:46:GLN:NE2	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ILE:HD11	1:A:162:PRO:HG3	1.86	0.57
1:B:80:ASN:OD1	1:B:80:ASN:C	2.47	0.57
3:D:2:BGC:C6	3:D:5:XYS:C1	2.82	0.56
1:B:207:CYS:CB	1:B:216:CYS:HG	2.18	0.55
1:B:233:ASP:H	1:B:236:GLN:NE2	2.05	0.54
1:B:104:ASN:HD22	1:B:105:VAL:H	1.55	0.53
1:A:210:SER:O	1:A:213:ALA:O	2.27	0.52
1:B:100:ILE:HD12	2:E:1:NAG:H82	1.92	0.51
1:A:100:ILE:HD12	2:C:1:NAG:H82	1.94	0.49
1:A:59:GLN:HG3	1:A:129:TYR:CE1	2.47	0.48
1:B:35:LEU:HG	1:B:43:THR:HG23	1.94	0.48
1:A:18:THR:HG21	1:A:46:GLN:HE21	1.77	0.48
1:A:237:TYR:CE1	1:A:270:ARG:HG3	2.49	0.47
3:F:2:BGC:O6	3:F:5:XYS:C2	2.56	0.47
1:A:85:GLU:OE2	1:A:87:ASP:OD1	2.33	0.46
1:B:155:LYS:NZ	4:B:308:HOH:O	2.47	0.46
1:A:182:ARG:HD3	1:A:186:GLU:HG2	2.00	0.44
1:B:79:GLN:NE2	4:B:301:HOH:O	2.09	0.44
1:B:227:LYS:O	1:B:227:LYS:HD3	2.17	0.44
1:A:141:PHE:CD1	1:A:149:ARG:HG2	2.54	0.43
1:A:233:ASP:H	1:A:236:GLN:HE21	1.65	0.43
1:B:85:GLU:OE2	1:B:87:ASP:OD1	2.36	0.42
1:A:263:PRO:HA	1:A:264:PRO:HD3	1.87	0.42
1:A:135:ASN:ND2	1:A:137:TYR:H	2.17	0.42
1:B:234:ALA:O	1:B:238:ARG:HG3	2.20	0.41
1:B:115:GLN:HE22	1:B:271:ASP:CG	2.29	0.41
1:B:155:LYS:HB3	4:B:341:HOH:O	2.21	0.40
1:A:182:ARG:HD3	1:A:186:GLU:CG	2.51	0.40
1:B:153:ASN:C	1:B:153:ASN:ND2	2.77	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	265/278 (95%)	256 (97%)	9 (3%)	0	100	100
1	B	263/278 (95%)	253 (96%)	10 (4%)	0	100	100
All	All	528/556 (95%)	509 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	235/241 (98%)	225 (96%)	10 (4%)	26	13
1	B	233/241 (97%)	224 (96%)	9 (4%)	28	16
All	All	468/482 (97%)	449 (96%)	19 (4%)	27	15

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	110	LYS
1	A	135	ASN
1	A	207	CYS
1	A	208	GLU
1	A	258	ARG
1	A	265	GLU
1	A	267	LYS
1	A	270	ARG
1	A	272	ILE
1	B	59	GLN
1	B	80	ASN
1	B	81	SER
1	B	82	GLU
1	B	153	ASN
1	B	227	LYS
1	B	267	LYS
1	B	270	ARG

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Mol	Chain	Res	Type
1	B	272	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	23	HIS
1	A	28	ASN
1	A	36	HIS
1	A	46	GLN
1	A	135	ASN
1	A	165	GLN
1	A	175	ASN
1	A	203	HIS
1	A	219	GLN
1	A	236	GLN
1	B	14	ASN
1	B	23	HIS
1	B	31	ASN
1	B	36	HIS
1	B	46	GLN
1	B	97	GLN
1	B	104	ASN
1	B	115	GLN
1	B	153	ASN
1	B	230	GLN
1	B	236	GLN
1	B	251	ASN

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 ⓘ

18 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	NAG	C	1	2,1	14,14,15	0.63	0	17,19,21	1.41	2 (11%)
2	NAG	C	2	2	14,14,15	0.72	0	17,19,21	1.01	1 (5%)
2	BMA	C	3	2	11,11,12	0.69	0	15,15,17	0.74	0
3	BGC	D	1	3	11,11,12	0.69	0	15,15,17	0.91	0
3	BGC	D	2	3	11,11,12	0.64	0	15,15,17	0.98	0
3	BGC	D	3	3	11,11,12	0.75	0	15,15,17	0.92	0
3	XYS	D	4	3	9,9,10	1.23	1 (11%)	10,12,14	0.92	1 (10%)
3	XYS	D	5	3	9,9,10	1.31	1 (11%)	10,12,14	1.24	1 (10%)
3	GAL	D	6	3	11,11,12	0.73	0	15,15,17	1.07	1 (6%)
2	NAG	E	1	2,1	14,14,15	0.65	0	17,19,21	1.30	2 (11%)
2	NAG	E	2	2	14,14,15	0.69	0	17,19,21	0.89	1 (5%)
2	BMA	E	3	2	11,11,12	0.66	0	15,15,17	0.89	0
3	BGC	F	1	3	11,11,12	0.77	0	15,15,17	0.95	0
3	BGC	F	2	3	11,11,12	0.67	0	15,15,17	0.70	0
3	BGC	F	3	3	11,11,12	0.90	1 (9%)	15,15,17	0.96	2 (13%)
3	XYS	F	4	3	9,9,10	1.17	1 (11%)	10,12,14	0.93	1 (10%)
3	XYS	F	5	3	9,9,10	1.08	1 (11%)	10,12,14	1.65	1 (10%)
3	GAL	F	6	3	11,11,12	0.63	0	15,15,17	1.01	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	NAG	C	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
3	BGC	D	1	3	-	0/2/19/22	0/1/1/1
3	BGC	D	2	3	-	0/2/19/22	0/1/1/1
3	BGC	D	3	3	-	0/2/19/22	0/1/1/1
3	XYS	D	4	3	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	XYS	D	5	3	-	-	0/1/1/1
3	GAL	D	6	3	-	0/2/19/22	0/1/1/1
2	NAG	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
3	BGC	F	1	3	-	0/2/19/22	0/1/1/1
3	BGC	F	2	3	-	0/2/19/22	0/1/1/1
3	BGC	F	3	3	-	0/2/19/22	0/1/1/1
3	XYS	F	4	3	-	-	0/1/1/1
3	XYS	F	5	3	-	-	0/1/1/1
3	GAL	F	6	3	-	0/2/19/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	5	XYS	O5-C1	-3.62	1.36	1.43
3	D	4	XYS	O5-C1	-3.29	1.36	1.43
3	F	4	XYS	O5-C1	-3.23	1.36	1.43
3	F	5	XYS	O5-C1	-2.97	1.37	1.43
3	F	3	BGC	O5-C1	-2.27	1.39	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	O5-C1-C2	-4.56	104.23	111.29
3	F	5	XYS	C5-O5-C1	3.69	117.37	111.42
3	D	5	XYS	C5-O5-C1	3.44	116.96	111.42
2	E	1	NAG	C1-C2-N2	2.96	115.09	110.43
2	C	2	NAG	O5-C1-C2	-2.91	106.80	111.29
3	D	6	GAL	C1-C2-C3	2.56	113.37	109.64
2	E	1	NAG	O5-C1-C2	-2.49	107.44	111.29
3	F	4	XYS	C5-O5-C1	2.46	115.39	111.42
3	F	3	BGC	C1-O5-C5	2.46	115.49	112.19
3	D	4	XYS	C5-O5-C1	2.26	115.07	111.42
2	E	2	NAG	O5-C1-C2	-2.25	107.81	111.29
2	C	1	NAG	O6-C6-C5	-2.09	104.21	111.33
3	F	3	BGC	O6-C6-C5	-2.06	104.31	111.33
3	F	6	GAL	C1-O5-C5	2.03	114.90	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

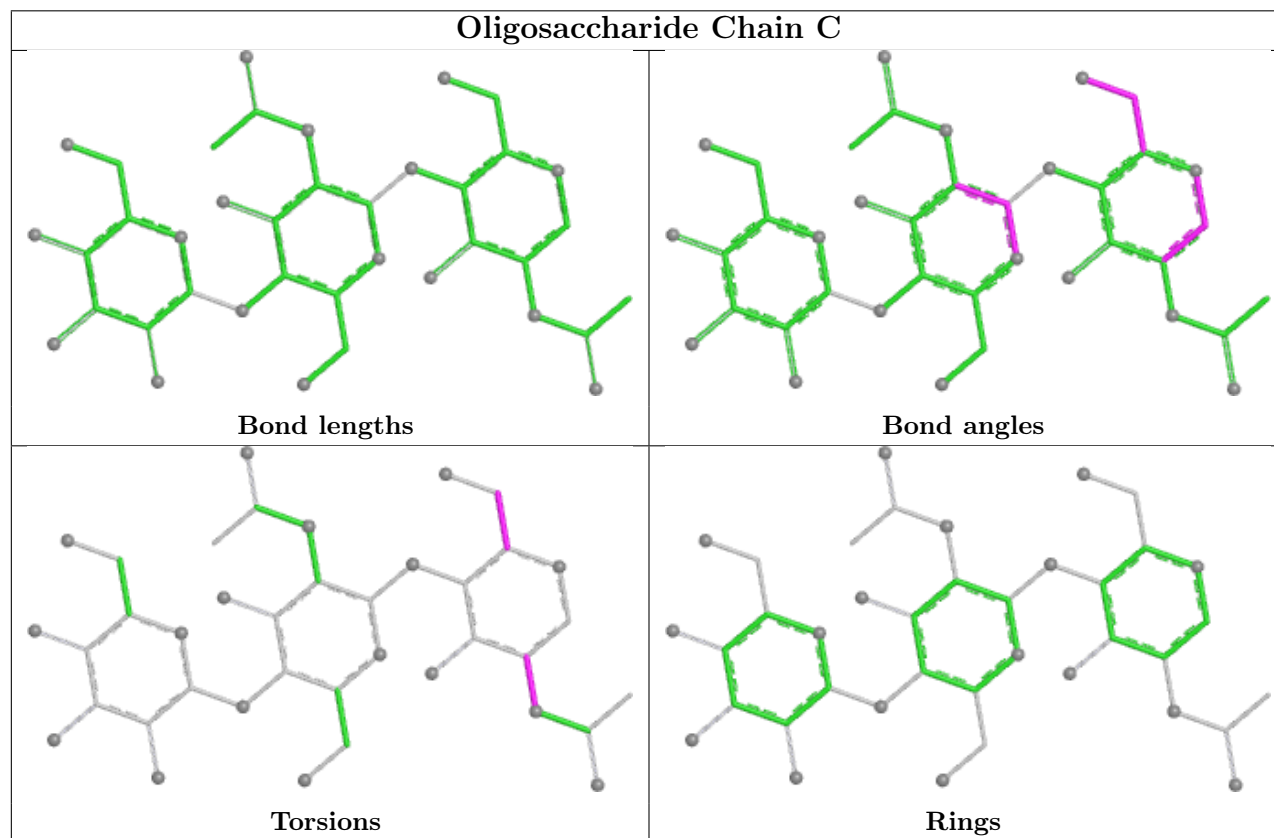
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C1-C2-N2-C7
2	E	1	NAG	C1-C2-N2-C7
2	C	1	NAG	C3-C2-N2-C7
2	E	1	NAG	C3-C2-N2-C7
2	C	1	NAG	C4-C5-C6-O6

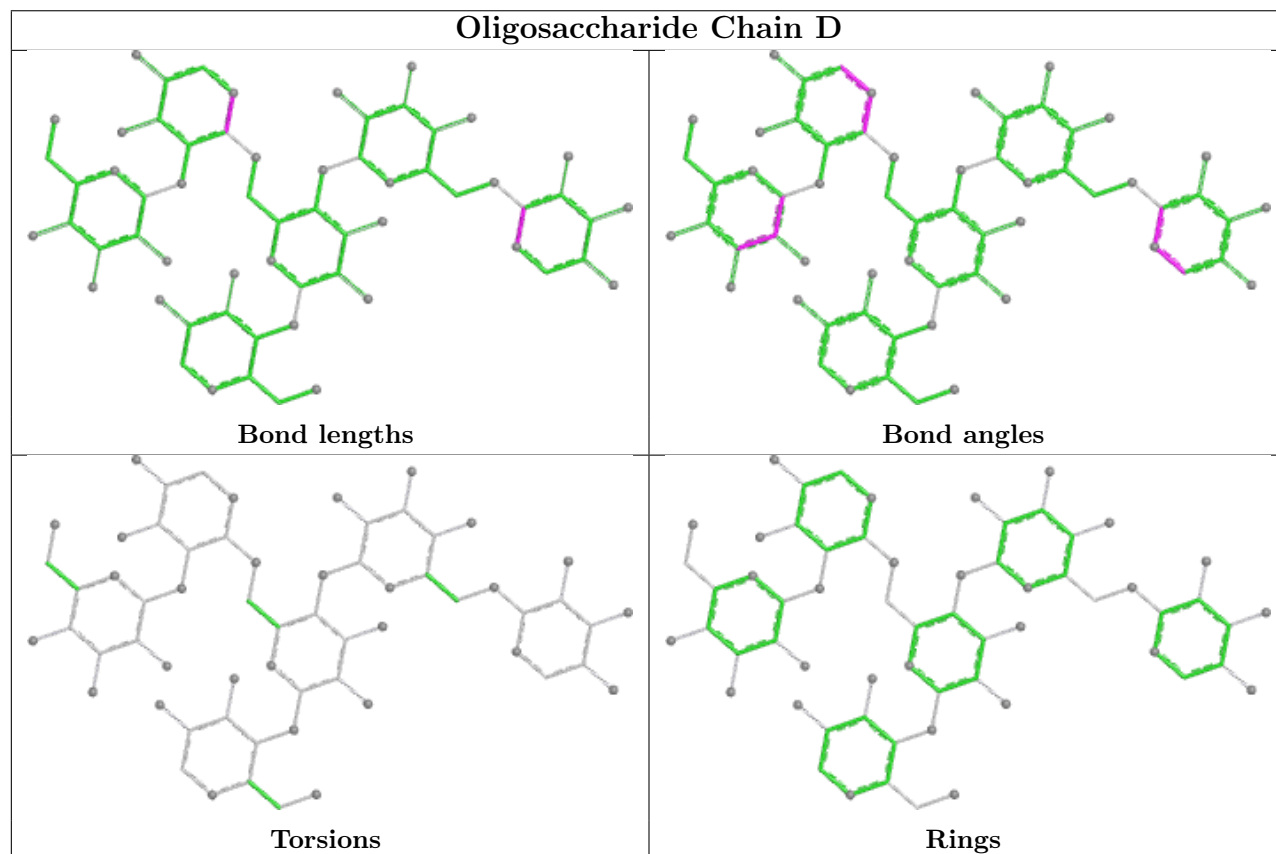
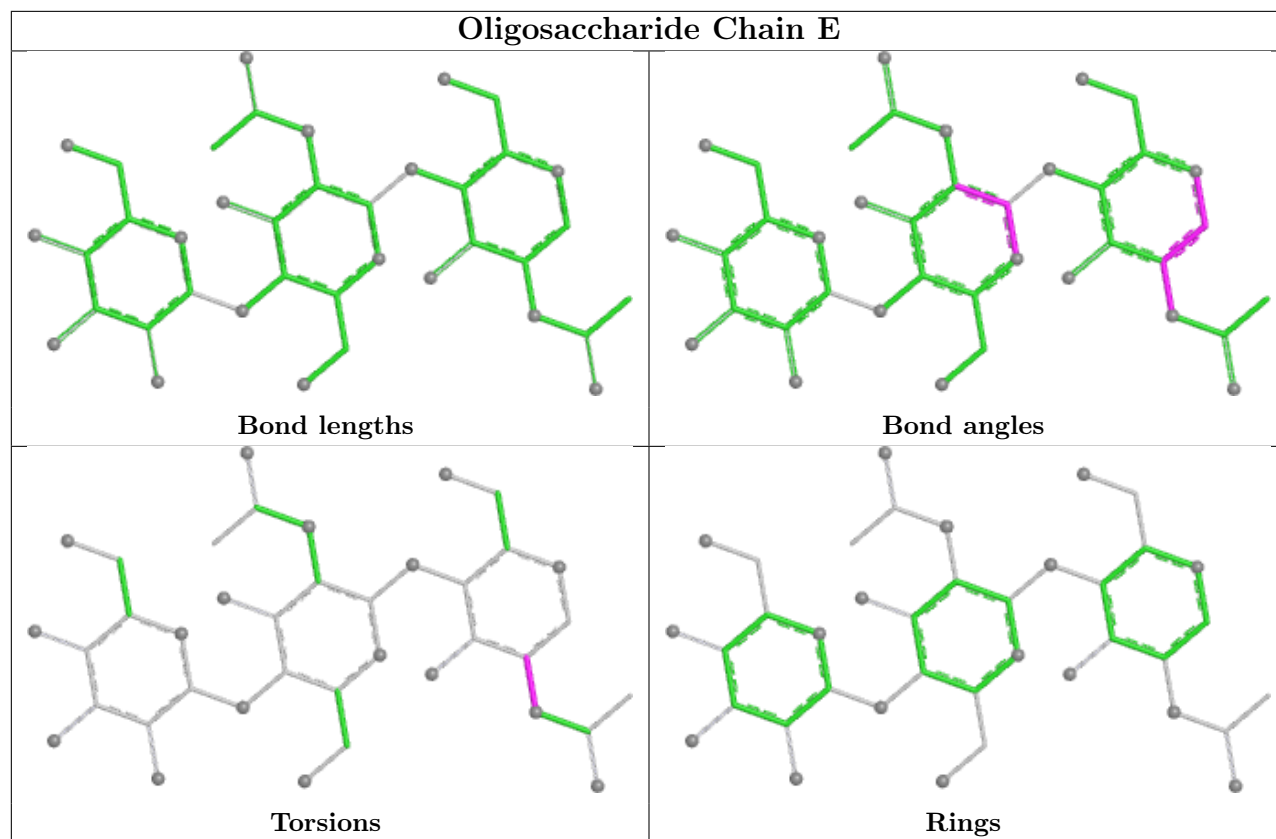
There are no ring outliers.

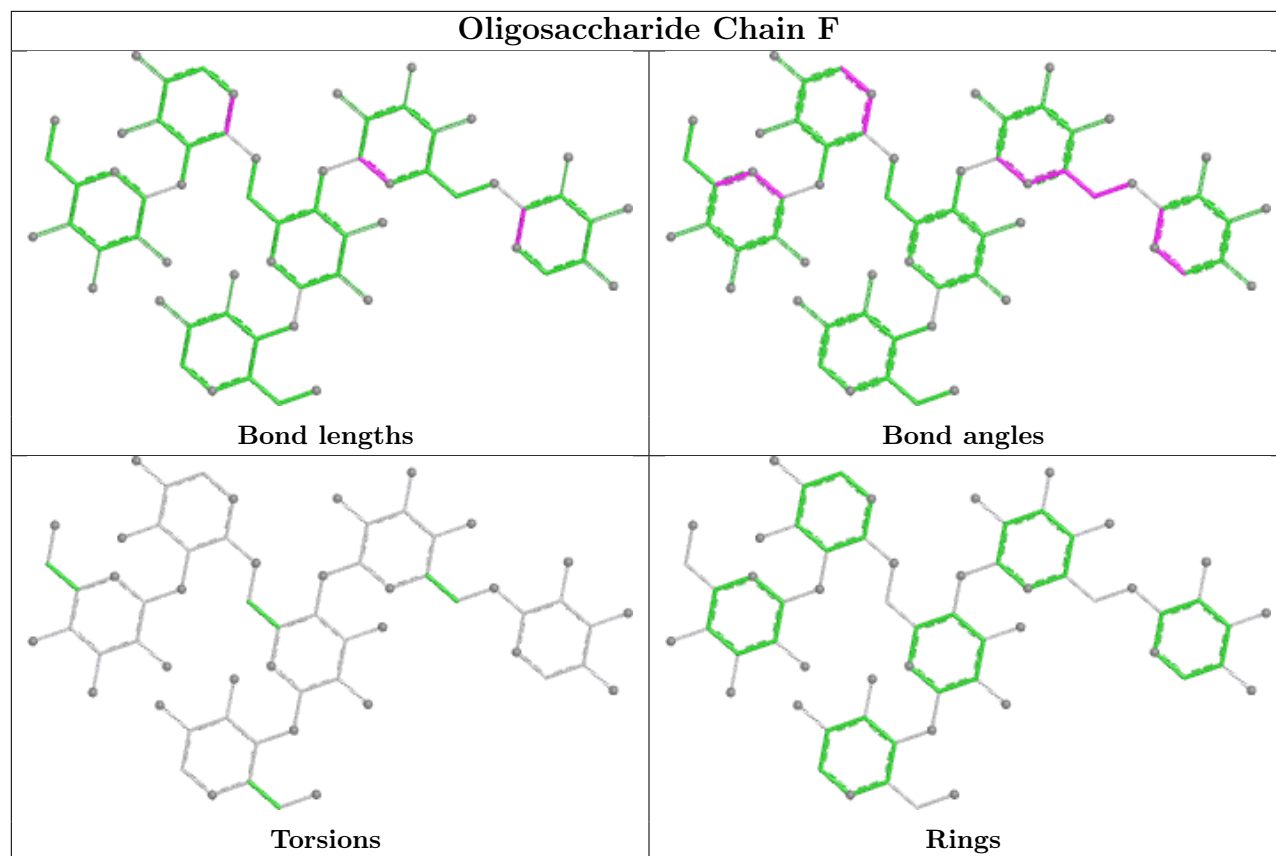
6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	1	NAG	3	0
3	D	5	XYS	2	0
2	C	1	NAG	1	0
3	F	2	BGC	2	0
3	F	5	XYS	2	0
3	D	2	BGC	2	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, 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	267/278 (96%)	0.91	49 (18%) 3 2	17, 29, 51, 70	4 (1%)
1	B	264/278 (94%)	0.56	25 (9%) 14 12	16, 28, 42, 49	3 (1%)
All	All	531/556 (95%)	0.74	74 (13%) 6 5	16, 29, 48, 70	7 (1%)

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	272	ILE	7.9
1	B	272	ILE	7.8
1	A	213	ALA	7.1
1	A	215	PHE	6.7
1	A	211	VAL	6.5
1	B	80	ASN	6.2
1	A	260	PRO	6.0
1	A	214	LYS	5.5
1	A	262	MET	5.0
1	A	253	CYS	4.8
1	B	9	VAL	4.7
1	A	259	TYR	4.7
1	A	209	ALA	4.6
1	A	261	SER	4.5
1	A	266	CYS	4.4
1	A	80	ASN	4.4
1	A	242	TRP	4.3
1	A	235	PHE	4.3
1	A	79	GLN	4.2
1	A	81	SER	4.1
1	B	81	SER	4.1
1	B	174[A]	TRP	4.0
1	A	263	PRO	3.8
1	A	210	SER	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	212	GLU	3.6
1	B	79	GLN	3.5
1	A	7	VAL	3.5
1	A	264	PRO	3.4
1	A	207	CYS	3.2
1	A	221	ALA	3.1
1	A	158	GLY	3.1
1	A	241	SER	3.1
1	B	10	ALA	3.0
1	A	220	GLY	3.0
1	A	216	CYS	2.9
1	A	258	ARG	2.9
1	A	267	LYS	2.8
1	A	178	ASP	2.8
1	B	156	ASP	2.8
1	A	28	ASN	2.7
1	A	6	PRO	2.7
1	B	137	TYR	2.6
1	B	27	PHE	2.6
1	A	208	GLU	2.6
1	B	145	ASP	2.6
1	A	218	THR	2.5
1	B	13	ARG	2.5
1	A	268	ARG	2.4
1	A	145	ASP	2.4
1	A	156	ASP	2.4
1	A	234	ALA	2.4
1	A	8	ASP	2.3
1	B	178	ASP	2.3
1	B	209	ALA	2.3
1	B	28	ASN	2.3
1	B	253	CYS	2.3
1	B	235	PHE	2.2
1	B	208	GLU	2.2
1	A	228	GLU	2.1
1	A	245	GLN	2.1
1	A	256	ARG	2.1
1	A	265	GLU	2.1
1	B	228	GLU	2.1
1	A	271	ASP	2.1
1	A	219	GLN	2.1
1	A	257	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	157	LEU	2.1
1	A	269	ASP	2.1
1	B	21	PHE	2.0
1	B	257	SER	2.0
1	B	259	TYR	2.0
1	A	243	VAL	2.0
1	B	261	SER	2.0
1	B	30	GLY	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.

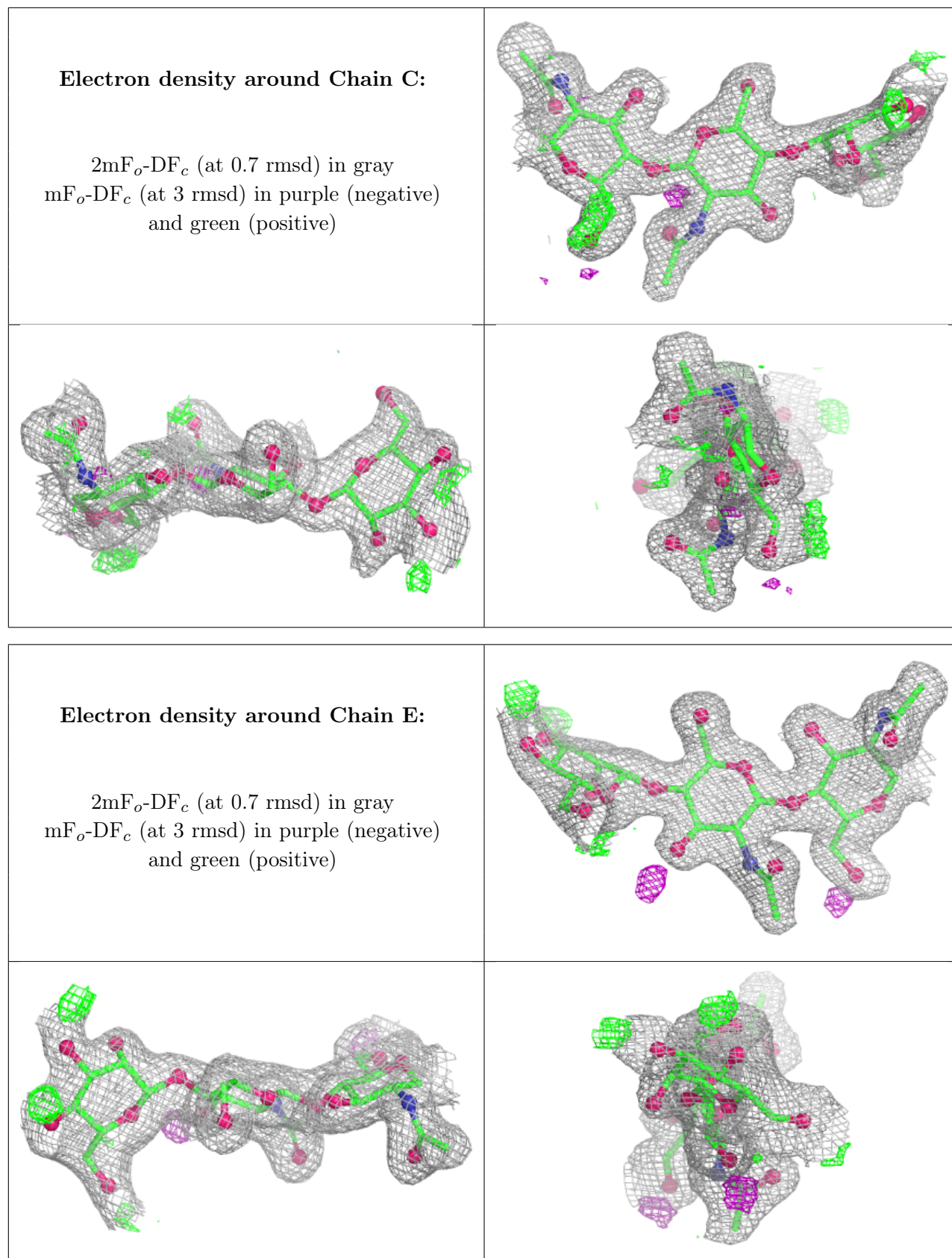
6.3 Carbohydrates [i](#)

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, 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($\text{\AA}^2$)	Q<0.9
3	BGC	D	1	11/12	0.64	0.19	48,50,52,54	0
2	BMA	E	3	11/12	0.75	0.15	48,52,54,57	0
2	BMA	C	3	11/12	0.77	0.15	51,54,57,59	0
3	GAL	D	6	11/12	0.80	0.17	54,55,56,56	0
3	GAL	F	6	11/12	0.86	0.12	39,42,43,45	0
3	XYS	D	4	9/10	0.87	0.11	47,48,48,49	0
3	BGC	F	1	11/12	0.87	0.12	34,38,43,45	0
3	XYS	D	5	9/10	0.87	0.13	44,47,49,50	0
3	XYS	F	4	9/10	0.90	0.10	38,40,43,45	0
2	NAG	C	2	14/15	0.90	0.11	35,38,41,45	0
2	NAG	C	1	14/15	0.92	0.09	23,26,32,32	0
2	NAG	E	2	14/15	0.93	0.09	30,34,38,41	0
3	BGC	D	3	11/12	0.93	0.09	29,35,39,42	0
3	XYS	F	5	9/10	0.94	0.09	25,32,35,37	0
3	BGC	D	2	11/12	0.94	0.09	38,41,43,44	0
2	NAG	E	1	14/15	0.96	0.07	22,26,31,32	0
3	BGC	F	2	11/12	0.96	0.07	26,29,32,34	0
3	BGC	F	3	11/12	0.96	0.06	22,27,29,31	0

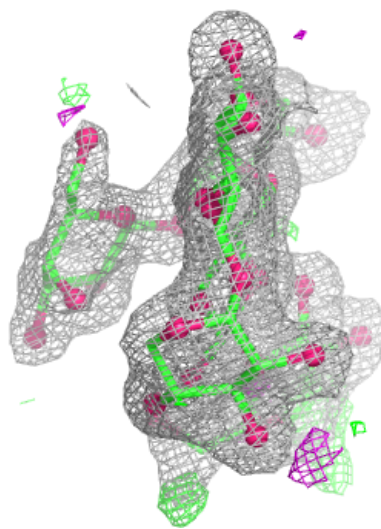
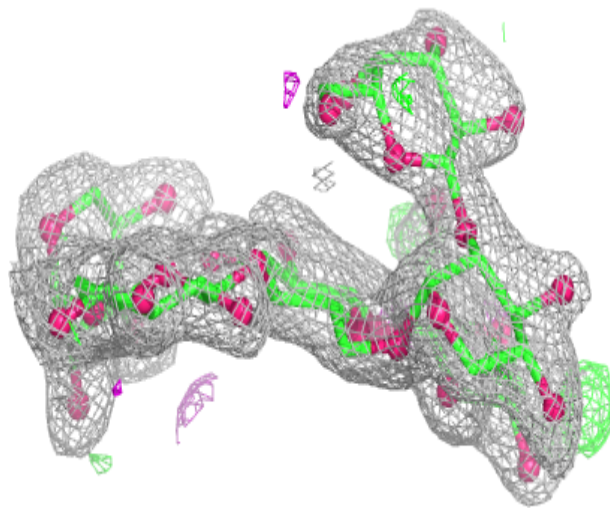
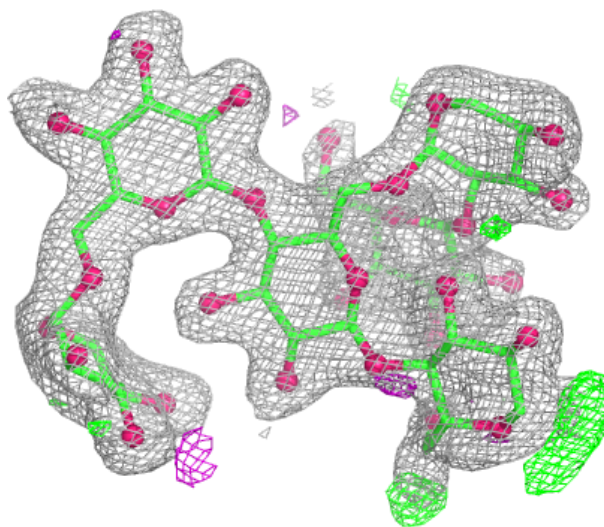
The following is a graphical depiction of the model fit to experimental electron density for oligosac-

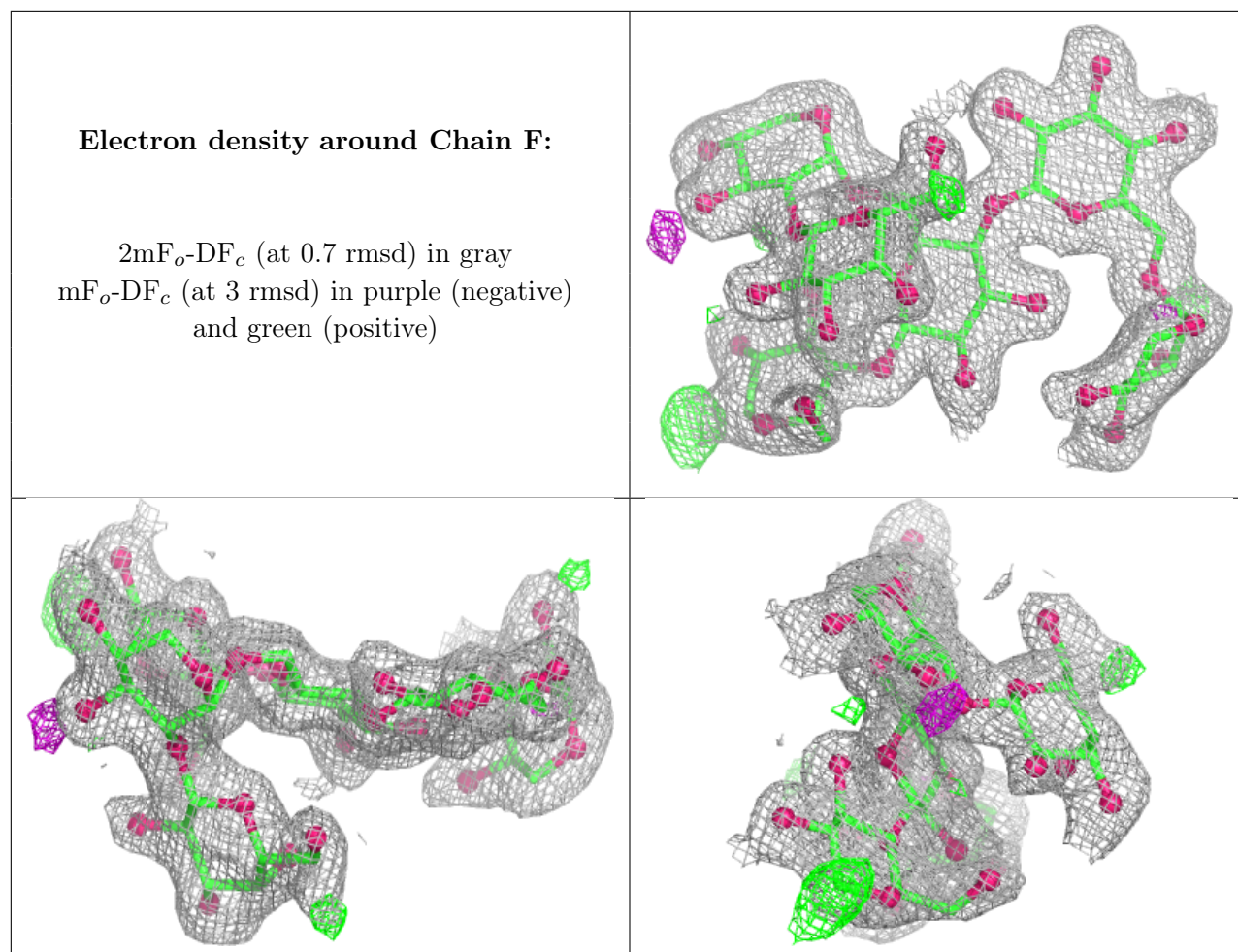
charide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain D:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.