



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:06 PM UTC

PDB ID : 5GM5 / pdb_00005gm5
Title : Crystal structure of FI-CMCase from *Aspergillus aculeatus* F-50 in complex with cellobiose
Authors : Huang, J.W.; Liu, W.D.; Zheng, Y.Y.; Chen, C.C.; Guo, R.T.
Deposited on : 2016-07-12
Resolution : 1.73 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

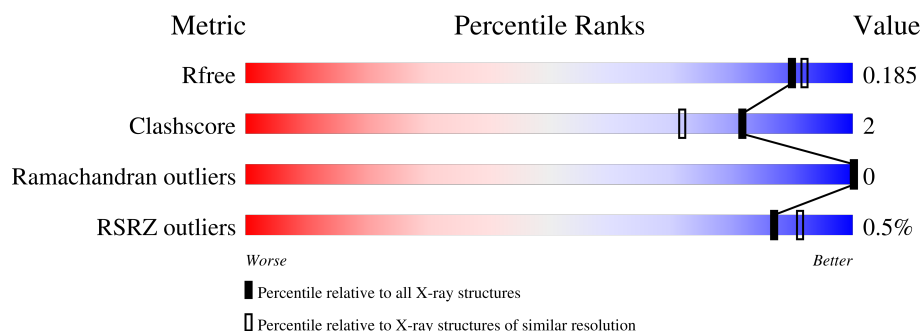
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

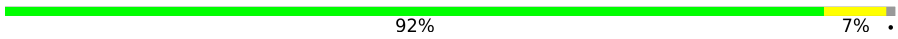
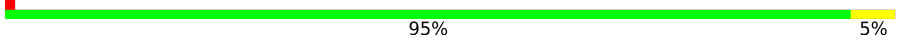
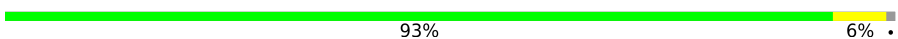
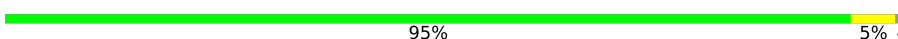
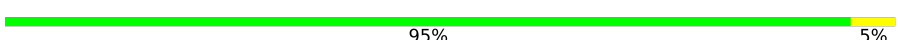
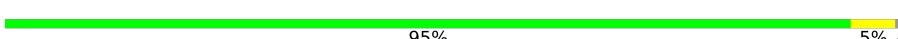
The reported resolution of this entry is 1.73 Å.

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	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	220	 92% 7% .
1	B	220	 95% 5% %
1	C	220	 93% 6% .
1	D	220	 95% 5% .
1	E	220	 95% 5%
1	F	220	 95% 5% .

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Mol	Chain	Length	Quality of chain
1	G	220	% 96% .
2	H	2	50% 50%
2	I	2	100%
2	J	2	50% 50%
2	K	2	100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13561 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Endoglucanase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	0	0
			1675	1062	276	334	3			
1	B	219	Total	C	N	O	S	0	0	0
			1680	1065	277	335	3			
1	C	218	Total	C	N	O	S	0	0	0
			1675	1062	276	334	3			
1	D	218	Total	C	N	O	S	0	0	0
			1675	1062	276	334	3			
1	E	220	Total	C	N	O	S	0	0	0
			1689	1070	279	337	3			
1	F	218	Total	C	N	O	S	0	0	0
			1675	1062	276	334	3			
1	G	219	Total	C	N	O	S	0	0	0
			1680	1065	277	335	3			

There are 7 discrepancies between the modelled and reference sequences:

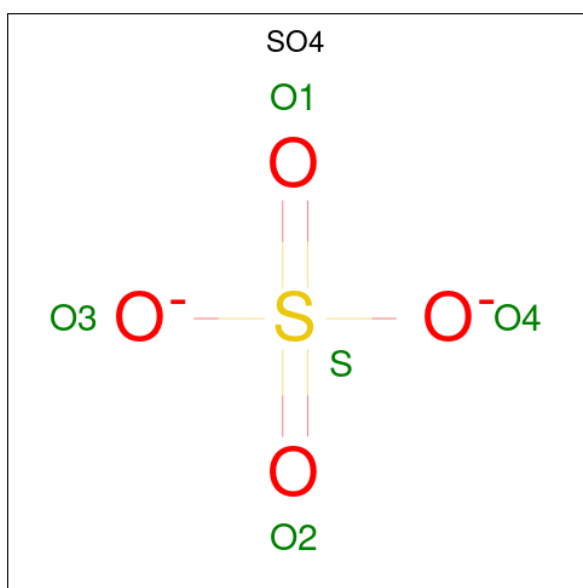
Chain	Residue	Modelled	Actual	Comment	Reference
A	202	ALA	GLU	engineered mutation	UNP P22669
B	202	ALA	GLU	engineered mutation	UNP P22669
C	202	ALA	GLU	engineered mutation	UNP P22669
D	202	ALA	GLU	engineered mutation	UNP P22669
E	202	ALA	GLU	engineered mutation	UNP P22669
F	202	ALA	GLU	engineered mutation	UNP P22669
G	202	ALA	GLU	engineered mutation	UNP P22669

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	H	2	Total	C	O	0	0	0
			23	12	11			
2	I	2	Total	C	O	0	0	0
			23	12	11			
2	J	2	Total	C	O	0	0	0
			23	12	11			
2	K	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



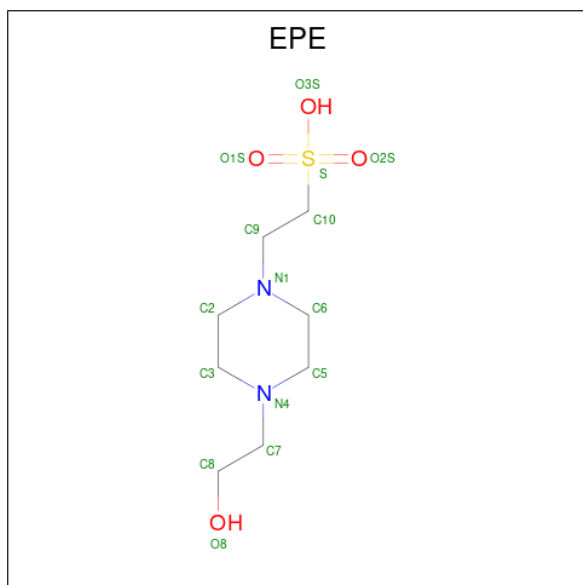
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	249	Total	O	0	0
			249	249		
5	B	243	Total	O	0	0
			243	243		
5	C	224	Total	O	0	0
			224	224		
5	D	269	Total	O	0	0
			269	269		
5	E	258	Total	O	0	0
			258	258		
5	F	205	Total	O	0	0
			205	205		
5	G	212	Total	O	0	0
			212	212		

3 Residue-property plots [i](#)

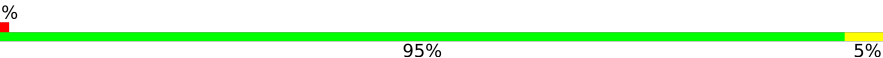
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

- Molecule 1: Endoglucanase-1

Chain A: 



- Molecule 1: Endoglucanase-1

Chain B: 

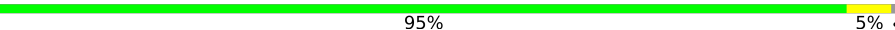


- Molecule 1: Endoglucanase-1

Chain C: 

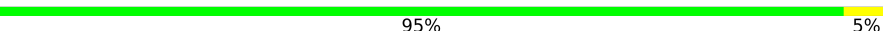


- Molecule 1: Endoglucanase-1

Chain D: 

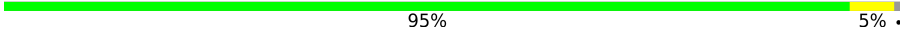


- Molecule 1: Endoglucanase-1

Chain E: 

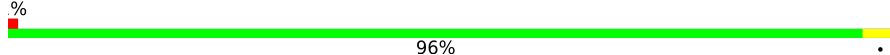


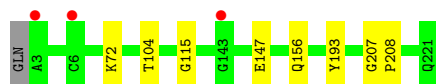
- Molecule 1: Endoglucanase-1

Chain F:  95% 5%



- Molecule 1: Endoglucanase-1

Chain G:  96%

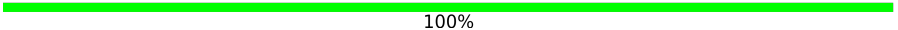


- Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain H:  50% 50%



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

Chain I:  100%

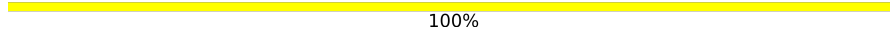


- Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain J:  50% 50%



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

Chain K:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.45Å 85.57Å 106.53Å 90.00° 91.87° 90.00°	Depositor
Resolution (Å)	25.00 – 1.73 25.00 – 1.73	Depositor EDS
% Data completeness (in resolution range)	99.7 (25.00-1.73) 99.7 (25.00-1.73)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 1.73Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.144 , 0.181 (Not available) , 0.185	Depositor DCC
R_{free} test set	8316 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for k,h,-l 0.005 for -k,-h,-l 0.019 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13561	wwPDB-VP
Average B, all atoms (Å ²)	21.0	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.*

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: BGC, SO4, EPE

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	0.71	0/1724	0.87	1/2359 (0.0%)
1	B	0.72	0/1729	0.89	0/2366
1	C	0.70	0/1724	0.83	0/2359
1	D	0.71	0/1724	0.86	0/2359
1	E	0.72	0/1738	0.90	1/2378 (0.0%)
1	F	0.63	0/1724	0.86	0/2359
1	G	0.66	0/1729	0.87	0/2366
All	All	0.69	0/12092	0.87	2/16546 (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	E	84	THR	CA-CB-OG1	5.16	117.34	109.60
1	A	38	ASN	N-CA-C	-5.02	106.85	113.12

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	1675	0	1556	11	0
1	B	1680	0	1561	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1675	0	1556	9	0
1	D	1675	0	1556	7	0
1	E	1689	0	1569	14	0
1	F	1675	0	1556	7	0
1	G	1680	0	1561	5	0
2	H	23	0	21	0	0
2	I	23	0	21	0	0
2	J	23	0	21	0	0
2	K	23	0	21	0	0
3	A	5	0	0	0	0
3	B	10	0	0	1	0
3	C	5	0	0	0	0
3	D	10	0	0	0	0
3	E	5	0	0	0	0
3	F	5	0	0	0	0
3	G	5	0	0	0	0
4	B	15	0	18	0	0
5	A	249	0	0	3	0
5	B	243	0	0	1	2
5	C	224	0	0	3	1
5	D	269	0	0	2	0
5	E	258	0	0	5	1
5	F	205	0	0	0	0
5	G	212	0	0	1	0
All	All	13561	0	11017	57	2

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 2.

All (57) 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:86:ARG:NH2	3:B:302:SO4:O2	2.16	0.77
1:C:147:GLU:OE1	5:C:401:HOH:O	2.04	0.74
1:A:7:ASP:OD1	5:A:401:HOH:O	2.07	0.72
1:C:86:ARG:NH1	5:C:402:HOH:O	2.23	0.70
1:G:147:GLU:OE2	5:G:401:HOH:O	2.12	0.68
1:E:157:LYS:HE2	5:E:546:HOH:O	1.96	0.65
1:D:147:GLU:OE2	5:D:401:HOH:O	2.14	0.63
1:C:115:GLY:H	1:C:156:GLN:HE21	1.44	0.63
1:E:119:LEU:HD23	1:E:119:LEU:C	2.24	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:145:THR:HG23	5:C:468:HOH:O	2.02	0.59
1:D:130:GLN:HE21	1:E:72:LYS:NZ	2.00	0.59
1:E:115:GLY:H	1:E:156:GLN:HE21	1.53	0.56
1:A:115:GLY:H	1:A:156:GLN:HE21	1.54	0.56
1:D:115:GLY:H	1:D:156:GLN:HE21	1.54	0.56
1:B:20:ASN:HB3	5:B:542:HOH:O	2.06	0.54
1:A:125:ARG:NH1	5:A:405:HOH:O	2.38	0.54
1:F:207:GLY:HA2	1:F:208:PRO:C	2.34	0.53
1:A:130:GLN:HE21	1:C:72:LYS:NZ	2.08	0.52
1:A:150:TYR:OH	1:A:157:LYS:HE3	2.09	0.52
1:B:207:GLY:HA2	1:B:208:PRO:C	2.34	0.52
1:C:207:GLY:HA2	1:C:208:PRO:C	2.35	0.51
1:D:4:GLN:N	5:D:407:HOH:O	2.44	0.51
1:E:84:THR:HG22	5:E:595:HOH:O	2.09	0.51
1:A:4:GLN:N	5:A:406:HOH:O	2.44	0.51
1:B:115:GLY:H	1:B:156:GLN:HE21	1.56	0.51
1:D:207:GLY:HA2	1:D:208:PRO:C	2.35	0.50
1:B:52:ASN:HD22	1:B:52:ASN:C	2.19	0.49
1:F:115:GLY:H	1:F:156:GLN:HE21	1.60	0.49
1:E:84:THR:HG22	1:E:171:GLN:HG3	1.94	0.49
1:F:130:GLN:HE21	1:G:72:LYS:NZ	2.10	0.49
1:E:84:THR:HG21	1:E:171:GLN:HE21	1.78	0.48
1:E:84:THR:CG2	5:E:595:HOH:O	2.60	0.48
1:C:80:GLN:HG3	1:C:82:PRO:HD3	1.96	0.47
1:A:19:ILE:HD12	1:A:19:ILE:N	2.30	0.47
1:B:153:ASN:C	1:B:153:ASN:OD1	2.58	0.46
1:G:115:GLY:H	1:G:156:GLN:HE21	1.63	0.46
1:E:119:LEU:C	1:E:119:LEU:CD2	2.89	0.44
1:E:207:GLY:HA2	1:E:208:PRO:C	2.43	0.44
1:E:119:LEU:HD23	1:E:120:MET:N	2.33	0.43
1:A:207:GLY:HA2	1:A:208:PRO:C	2.43	0.43
1:F:4:GLN:HE21	1:F:34:CYS:HB3	1.83	0.43
1:G:104:THR:HA	1:G:193:TYR:O	2.19	0.43
1:B:101:ASP:OD1	1:B:118:GLU:OE2	2.37	0.42
1:C:153:ASN:OD1	1:C:153:ASN:C	2.62	0.42
1:D:101:ASP:OD1	1:D:118:GLU:OE2	2.37	0.42
1:C:101:ASP:OD1	1:C:118:GLU:OE2	2.38	0.42
1:F:19:ILE:HD12	1:F:19:ILE:N	2.34	0.42
1:A:111:VAL:HG12	1:B:135:GLN:HE22	1.85	0.41
1:E:153:ASN:OD1	1:E:153:ASN:C	2.63	0.41
1:E:157:LYS:HD2	5:E:403:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:73:LYS:HG3	1:F:78:ILE:HG12	2.03	0.41
1:A:101:ASP:OD1	1:A:118:GLU:OE2	2.38	0.41
1:D:153:ASN:OD1	1:D:153:ASN:C	2.63	0.41
1:E:157:LYS:CE	5:E:546:HOH:O	2.60	0.41
1:G:207:GLY:HA2	1:G:208:PRO:C	2.46	0.41
1:A:153:ASN:OD1	1:A:153:ASN:C	2.62	0.40
1:F:207:GLY:CA	1:F:208:PRO:C	2.94	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:523:HOH:O	5:C:545:HOH:O[1_655]	2.19	0.01
5:B:585:HOH:O	5:E:414:HOH:O[1_655]	2.19	0.01

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	216/220 (98%)	213 (99%)	3 (1%)	0	100	100
1	B	217/220 (99%)	214 (99%)	3 (1%)	0	100	100
1	C	216/220 (98%)	213 (99%)	3 (1%)	0	100	100
1	D	216/220 (98%)	212 (98%)	4 (2%)	0	100	100
1	E	218/220 (99%)	212 (97%)	6 (3%)	0	100	100
1	F	216/220 (98%)	213 (99%)	3 (1%)	0	100	100
1	G	217/220 (99%)	214 (99%)	3 (1%)	0	100	100
All	All	1516/1540 (98%)	1491 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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	H	1	2	12,12,12	0.69	0	17,17,17	0.66	0
2	BGC	H	2	2	11,11,12	0.58	0	15,15,17	1.44	2 (13%)
2	BGC	I	1	2	12,12,12	0.56	0	17,17,17	0.61	0
2	BGC	I	2	2	11,11,12	0.43	0	15,15,17	0.96	0
2	BGC	J	1	2	12,12,12	0.60	0	17,17,17	0.72	0
2	BGC	J	2	2	11,11,12	0.50	0	15,15,17	1.60	3 (20%)
2	BGC	K	1	2	12,12,12	0.66	0	17,17,17	1.12	1 (5%)
2	BGC	K	2	2	11,11,12	0.57	0	15,15,17	1.37	3 (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	BGC	H	1	2	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	H	2	2	-	0/2/19/22	0/1/1/1
2	BGC	I	1	2	-	0/2/22/22	0/1/1/1
2	BGC	I	2	2	-	0/2/19/22	0/1/1/1
2	BGC	J	1	2	-	0/2/22/22	0/1/1/1
2	BGC	J	2	2	-	0/2/19/22	0/1/1/1
2	BGC	K	1	2	-	0/2/22/22	0/1/1/1
2	BGC	K	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	BGC	O2-C2-C3	-3.26	103.41	110.15
2	H	2	BGC	O2-C2-C3	-3.02	103.90	110.15
2	K	2	BGC	C1-O5-C5	2.97	116.17	112.19
2	J	2	BGC	C1-O5-C5	2.92	116.10	112.19
2	H	2	BGC	C1-O5-C5	2.90	116.07	112.19
2	J	2	BGC	C1-C2-C3	2.86	113.81	109.64
2	K	2	BGC	C6-C5-C4	-2.54	106.79	113.02
2	K	2	BGC	O4-C4-C5	-2.19	103.92	109.32
2	K	1	BGC	O5-C1-C2	2.02	113.85	110.30

There are no chirality outliers.

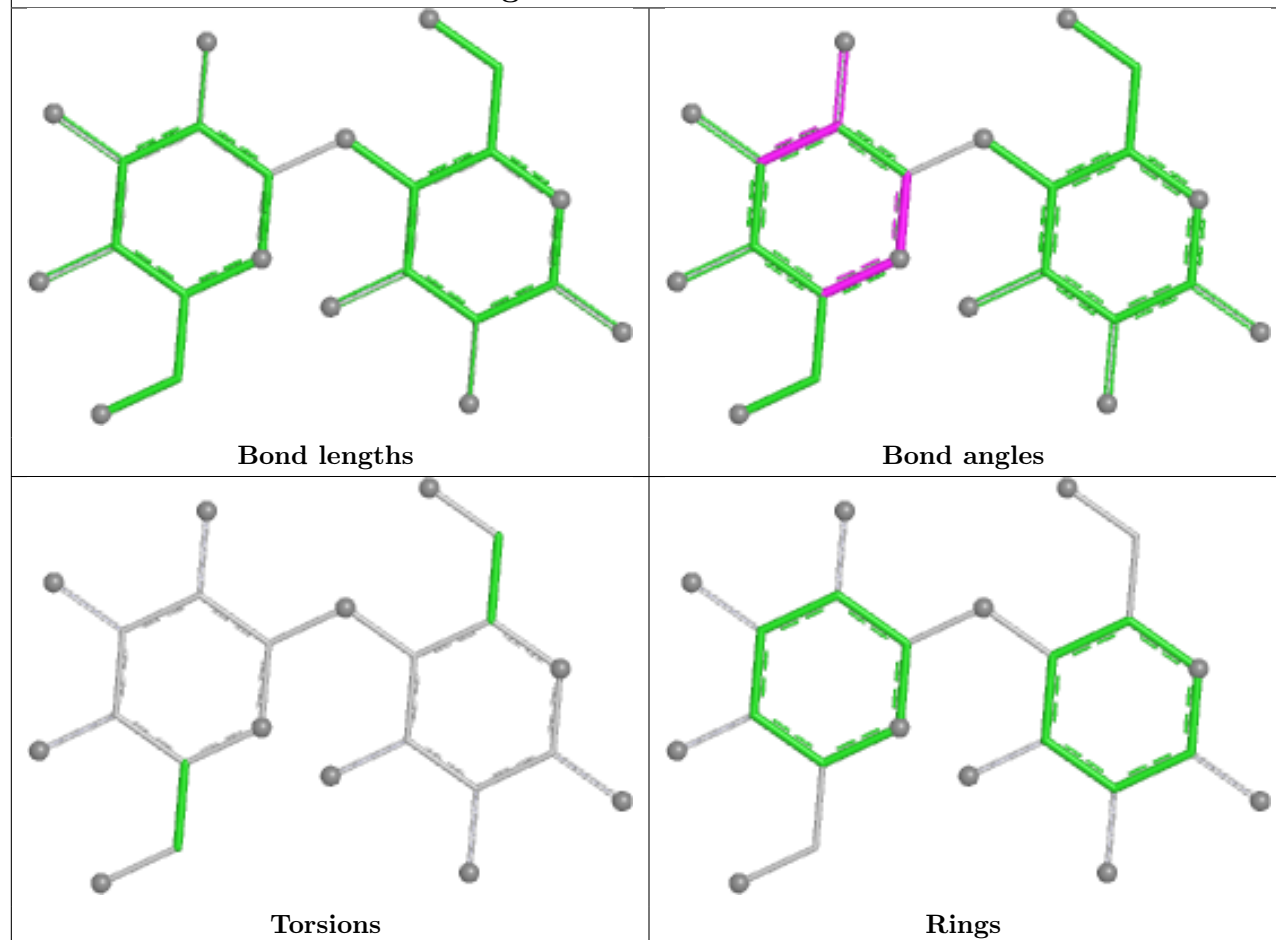
There are no torsion outliers.

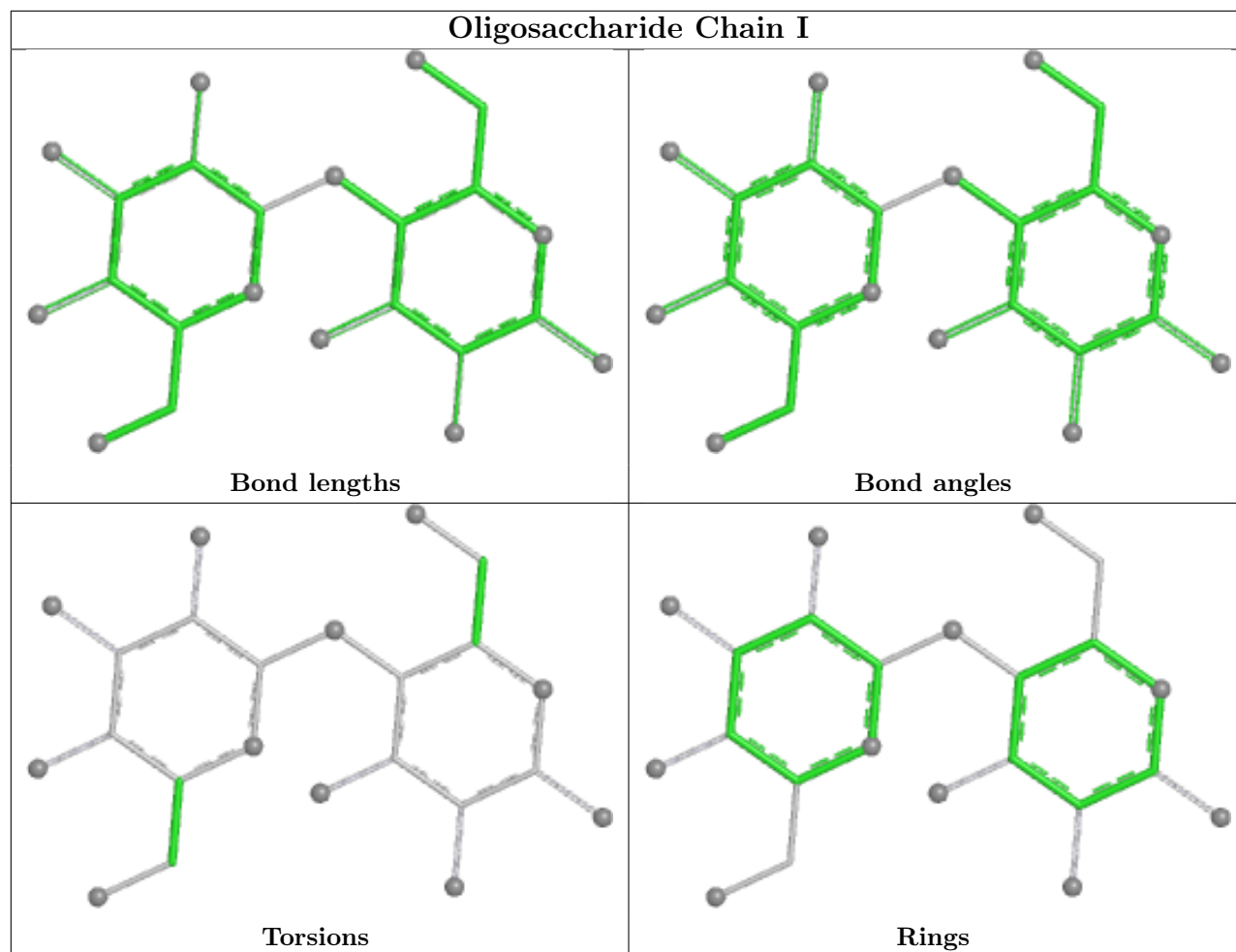
There are no ring outliers.

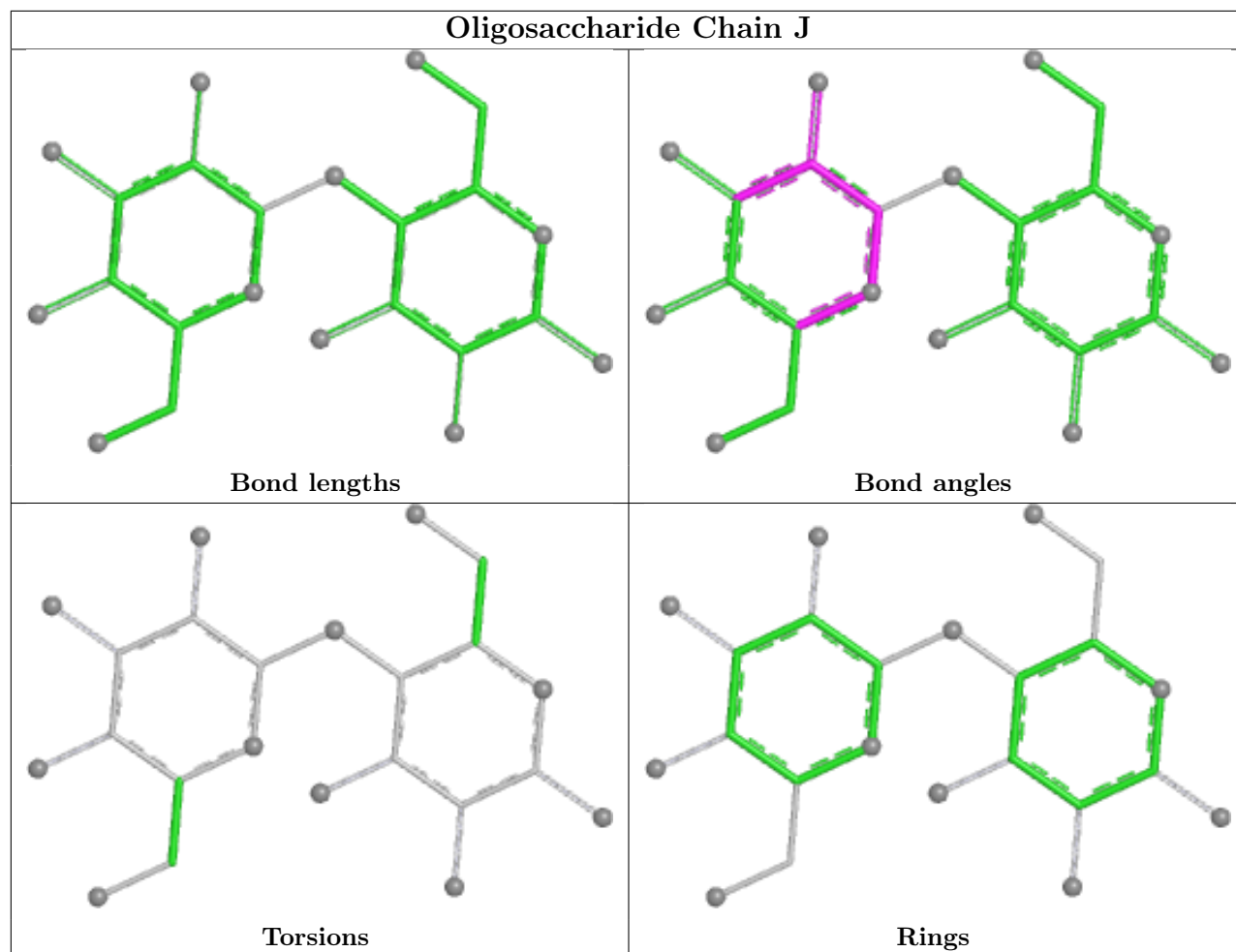
No monomer is involved in short contacts.

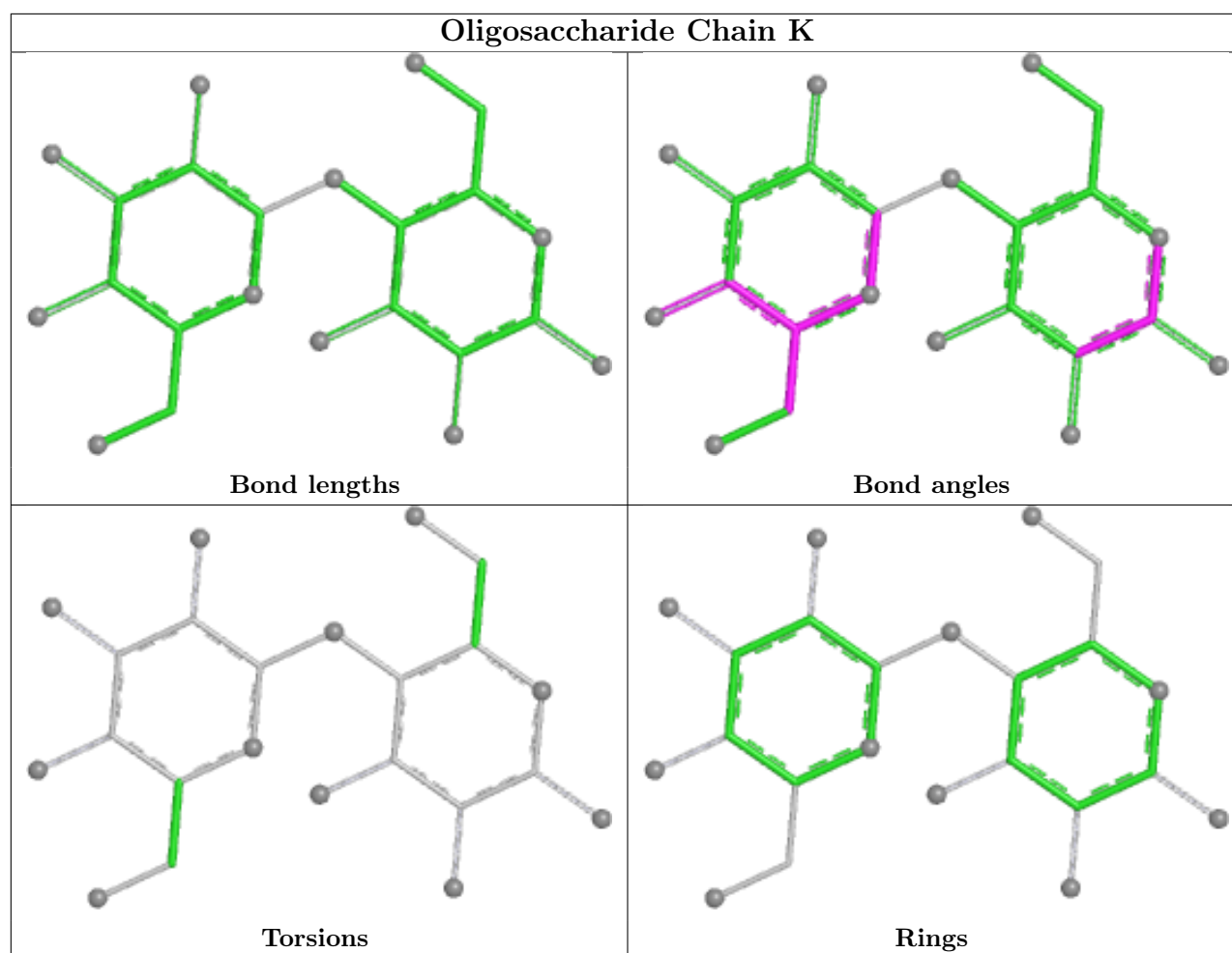
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain H









5.6 Ligand geometry [i](#)

10 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	SO4	B	301	-	4,4,4	0.56	0	6,6,6	0.30	0
3	SO4	D	303	-	4,4,4	0.42	0	6,6,6	0.32	0
4	EPE	B	303	-	15,15,15	1.81	1 (6%)	19,20,20	1.48	4 (21%)
3	SO4	F	302	-	4,4,4	0.42	0	6,6,6	0.24	0
3	SO4	A	302	-	4,4,4	0.55	0	6,6,6	0.35	0
3	SO4	B	302	-	4,4,4	0.47	0	6,6,6	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	C	301	-	4,4,4	0.48	0	6,6,6	0.75	0
3	SO4	E	301	-	4,4,4	0.49	0	6,6,6	0.21	0
3	SO4	G	302	-	4,4,4	0.52	0	6,6,6	0.21	0
3	SO4	D	302	-	4,4,4	0.56	0	6,6,6	0.43	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
4	EPE	B	303	-	-	3/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	303	EPE	C10-S	-6.57	1.68	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	303	EPE	O3S-S-C10	3.34	112.55	106.00
4	B	303	EPE	O2S-S-C10	-2.97	102.25	106.73
4	B	303	EPE	C7-N4-C3	-2.58	104.36	111.24
4	B	303	EPE	C3-C2-N1	-2.23	106.16	110.65

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	303	EPE	C9-C10-S-O2S
4	B	303	EPE	C9-C10-S-O3S
4	B	303	EPE	C9-C10-S-O1S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	SO4	1	0

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, 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	218/220 (99%)	-0.49	0 100 100	11, 17, 30, 39	0
1	B	219/220 (99%)	-0.44	2 (0%) 81 87	12, 17, 32, 44	0
1	C	218/220 (99%)	-0.40	1 (0%) 87 91	12, 18, 32, 44	0
1	D	218/220 (99%)	-0.54	0 100 100	11, 16, 27, 40	0
1	E	220/220 (100%)	-0.47	1 (0%) 87 91	11, 17, 27, 52	0
1	F	218/220 (99%)	-0.22	1 (0%) 87 91	15, 22, 37, 48	0
1	G	219/220 (99%)	-0.23	3 (1%) 73 80	13, 20, 33, 56	0
All	All	1530/1540 (99%)	-0.40	8 (0%) 87 91	11, 18, 32, 56	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	3	ALA	4.0
1	B	141	VAL	3.0
1	F	143	GLY	2.4
1	G	143	GLY	2.3
1	G	6	CYS	2.2
1	E	113	TRP	2.2
1	B	34	CYS	2.0
1	C	4	GLN	2.0

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

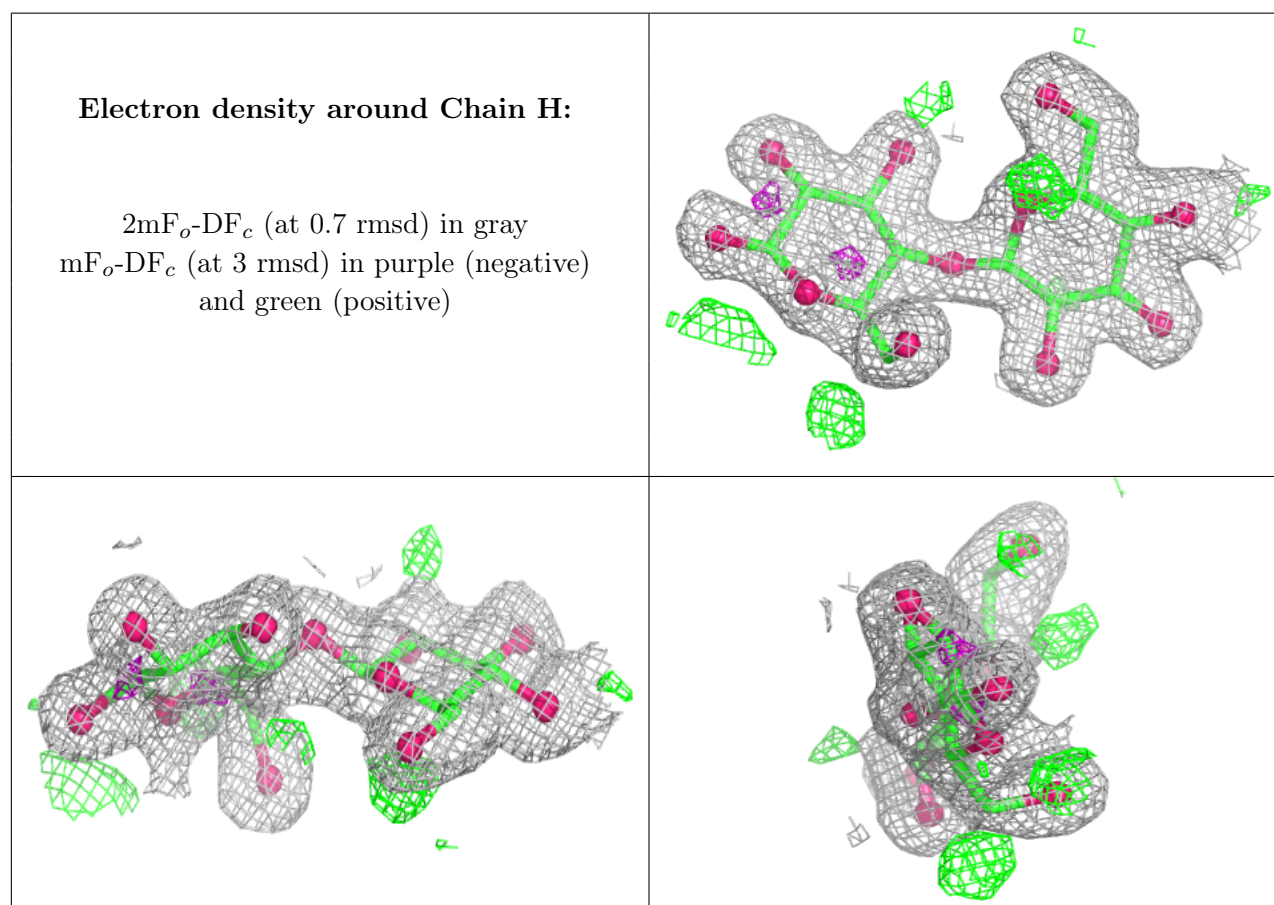
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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.

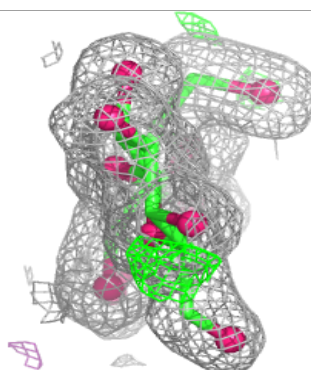
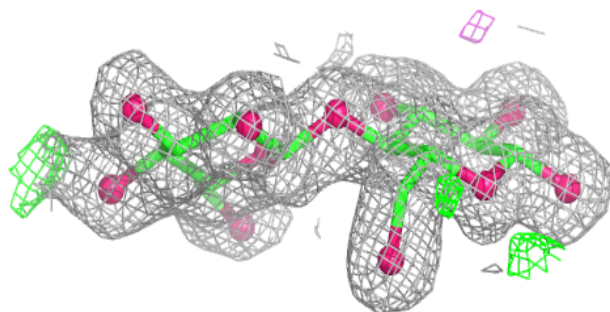
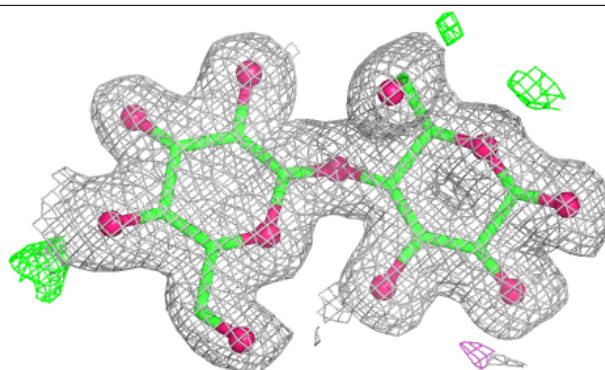
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BGC	J	1	12/12	0.87	0.10	28,35,39,43	0
2	BGC	J	2	11/12	0.89	0.09	29,35,42,45	0
2	BGC	H	2	11/12	0.94	0.07	20,24,27,30	0
2	BGC	K	2	11/12	0.94	0.07	21,26,28,31	0
2	BGC	H	1	12/12	0.95	0.07	18,26,30,31	0
2	BGC	I	2	11/12	0.96	0.06	19,22,29,33	0
2	BGC	K	1	12/12	0.96	0.05	19,21,24,26	0
2	BGC	I	1	12/12	0.96	0.05	17,20,22,24	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.

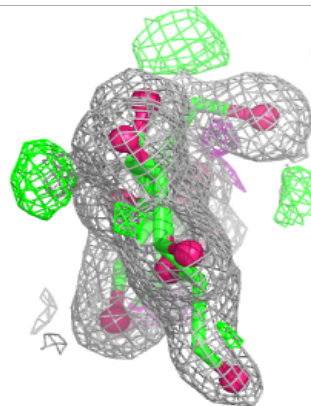
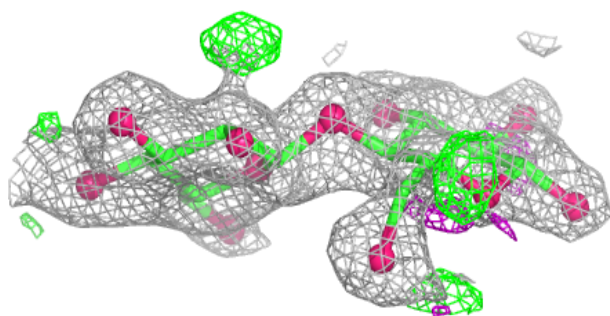
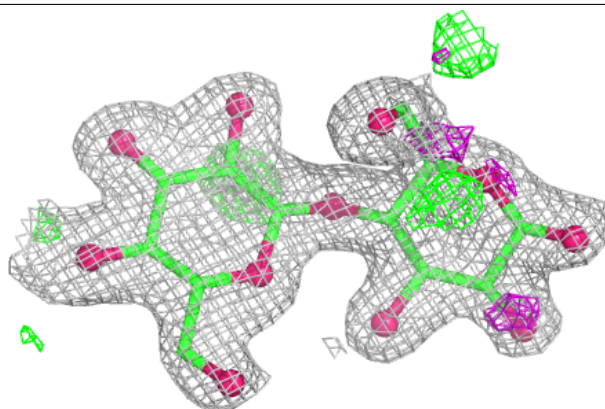


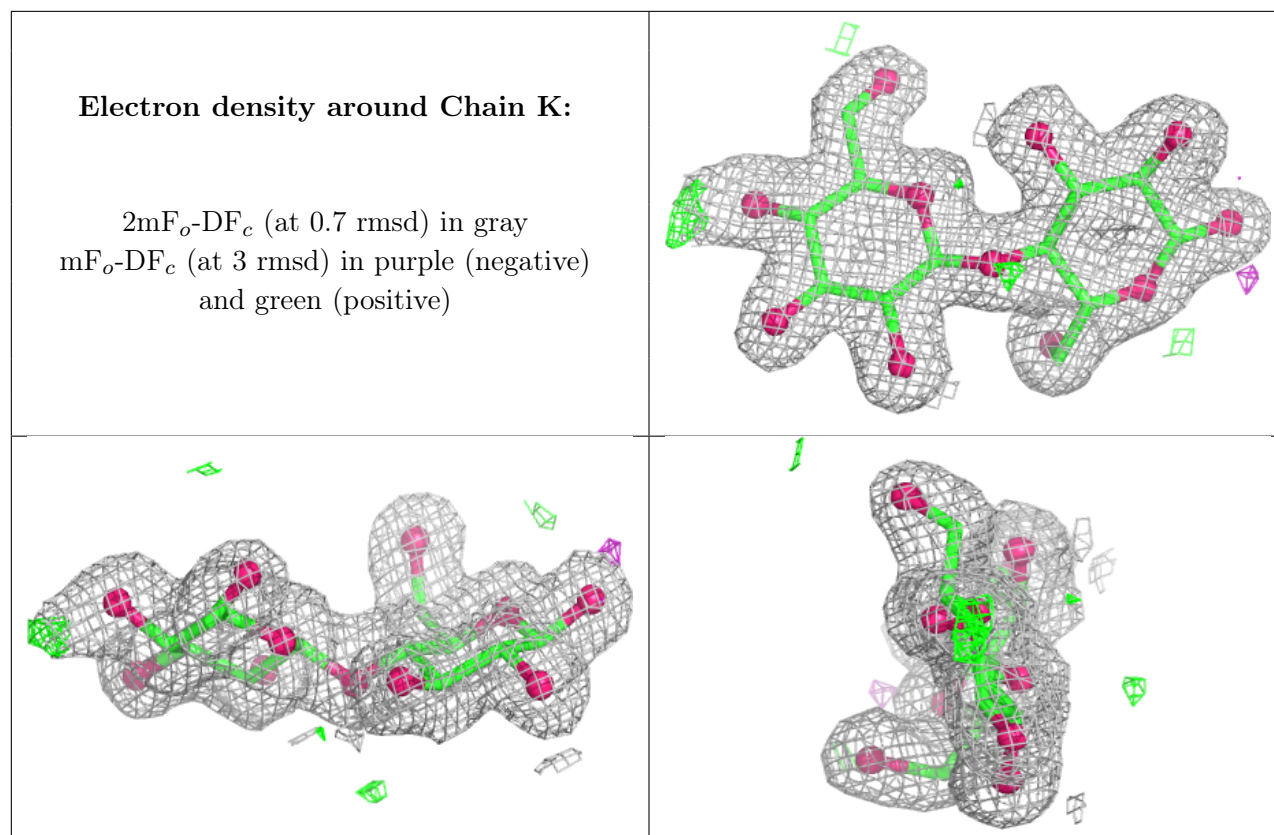
Electron density around Chain I:

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

**Electron density around Chain J:**

$2mF_o - DF_c$ (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($\text{\AA}^2$)	Q<0.9
3	SO4	G	302	5/5	0.75	0.10	54,63,83,95	0
3	SO4	C	301	5/5	0.78	0.13	34,50,75,81	0
3	SO4	E	301	5/5	0.83	0.14	46,64,71,83	0
3	SO4	B	302	5/5	0.83	0.12	56,61,65,86	0
3	SO4	D	303	5/5	0.84	0.10	51,63,76,90	0
4	EPE	B	303	15/15	0.84	0.14	35,47,53,57	0
3	SO4	B	301	5/5	0.86	0.09	43,46,60,74	0
3	SO4	A	302	5/5	0.89	0.11	32,53,63,64	0
3	SO4	D	302	5/5	0.89	0.10	32,52,65,69	0
3	SO4	F	302	5/5	0.93	0.08	56,58,72,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.