



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 12:34 PM UTC

PDB ID : 3EQO / pdb_00003eqo
Title : Crystal structure of beta-1,3-glucanase from *Phanerochaete chrysosporium* (Lam55A) gluconolactone complex
Authors : Ishida, T.; Fushinobu, S.; Kawai, R.; Kitaoka, M.; Igarashi, K.; Samejima, M.
Deposited on : 2008-10-01
Resolution : 2.25 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

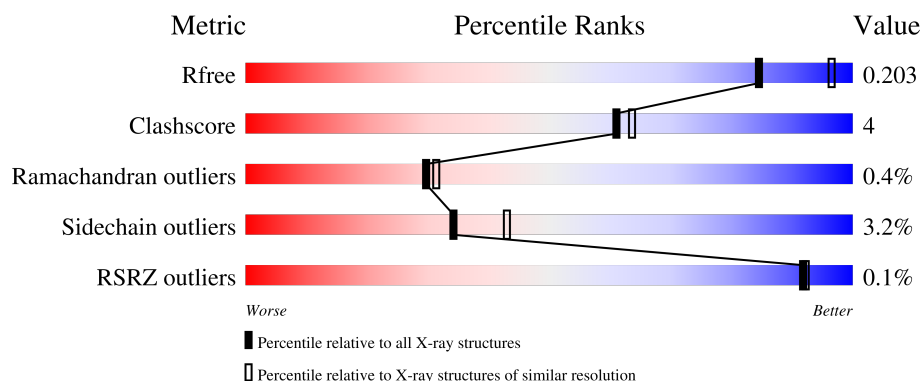
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	758	86% 12% ..
1	B	758	86% 12% ..
2	C	3	100%
2	D	3	100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12519 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.

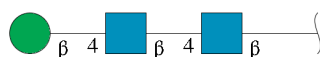
- Molecule 1 is a protein called Glucan 1,3-beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	752	Total	C	N	O	S	0	0	0
			5611	3526	973	1096	16			
1	B	752	Total	C	N	O	S	0	0	0
			5611	3526	973	1096	16			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLU	-	expression tag	UNP Q2Z1W1
A	-4	ALA	-	expression tag	UNP Q2Z1W1
A	-3	GLU	-	expression tag	UNP Q2Z1W1
A	-2	ALA	-	expression tag	UNP Q2Z1W1
A	-1	GLU	-	expression tag	UNP Q2Z1W1
A	0	PHE	-	expression tag	UNP Q2Z1W1
B	-5	GLU	-	expression tag	UNP Q2Z1W1
B	-4	ALA	-	expression tag	UNP Q2Z1W1
B	-3	GLU	-	expression tag	UNP Q2Z1W1
B	-2	ALA	-	expression tag	UNP Q2Z1W1
B	-1	GLU	-	expression tag	UNP Q2Z1W1
B	0	PHE	-	expression tag	UNP Q2Z1W1

- 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			

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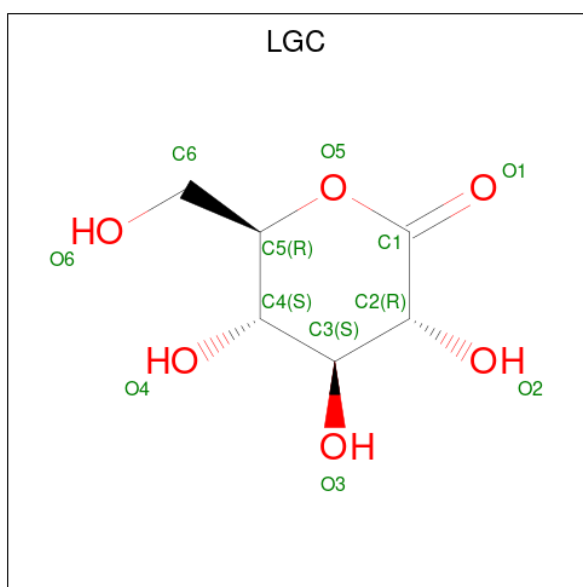
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is D-glucono-1,5-lactone (CCD ID: LGC) (formula: C₆H₁₀O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		
4	B	1	Total	C	O	0	0
			12	6	6		

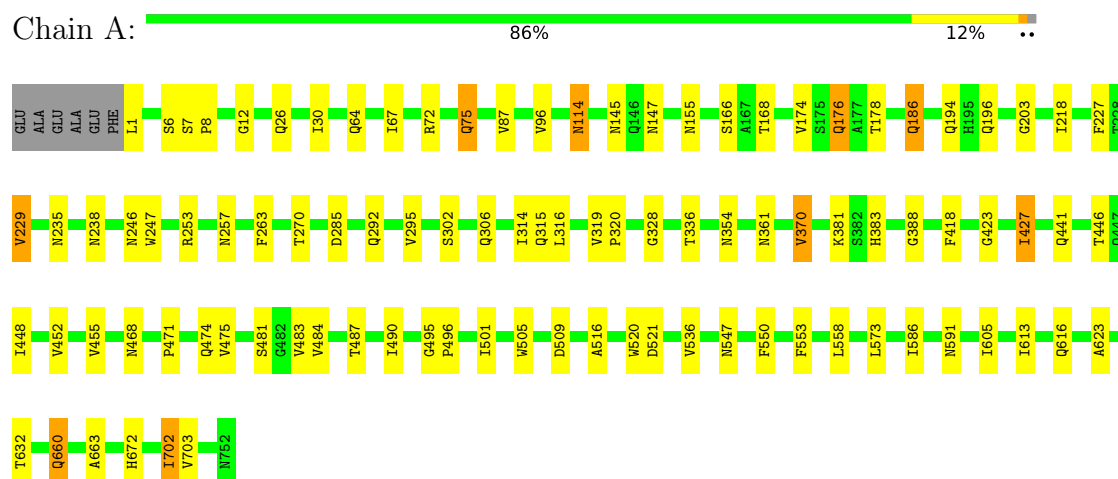
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	609	Total	O	0	0
			609	609		
5	B	584	Total	O	0	0
			584	584		

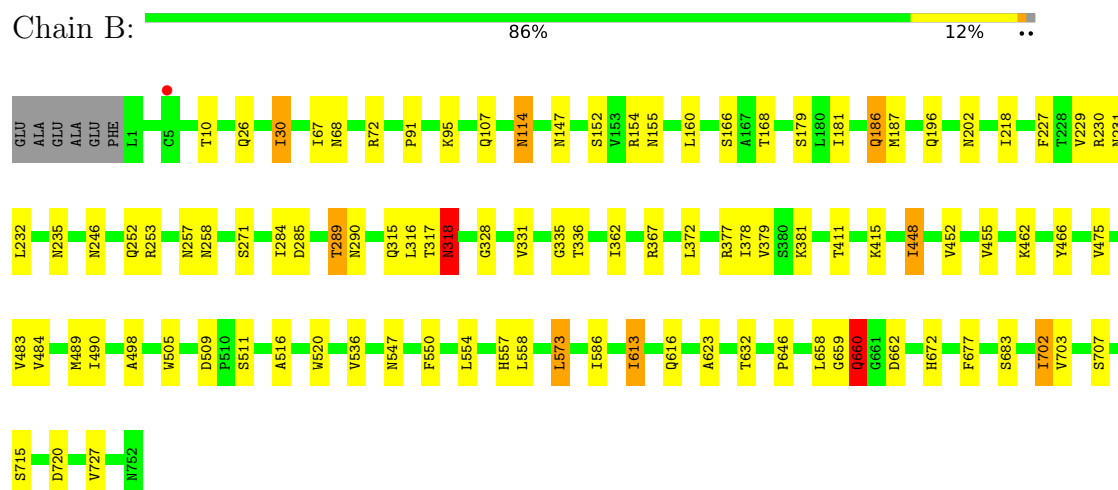
3 Residue-property plots

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: Glucan 1,3-beta-glucosidase



- Molecule 1: Glucan 1,3-beta-glucosidase



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



MAG1
MAG2
BMA3

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

Chain D:

100%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.53Å 67.15Å 109.78Å 93.94° 106.76° 97.08°	Depositor
Resolution (Å)	31.50 – 2.25 31.50 – 2.25	Depositor EDS
% Data completeness (in resolution range)	91.5 (31.50-2.25) 91.5 (31.50-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.26Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.142 , 0.200 0.144 , 0.203	Depositor DCC
R_{free} test set	3875 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.5	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12519	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 4.63% 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: ZN, LGC, NAG, BMA

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	1.21	6/5755 (0.1%)	1.11	13/7869 (0.2%)
1	B	1.18	6/5755 (0.1%)	1.08	7/7869 (0.1%)
All	All	1.20	12/11510 (0.1%)	1.09	20/15738 (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	B	0	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	452	VAL	CA-CB	7.36	1.64	1.54
1	B	613	ILE	CA-CB	6.53	1.62	1.54
1	A	229	VAL	C-O	-6.12	1.17	1.24
1	B	379	VAL	CA-CB	6.09	1.61	1.54
1	A	446	THR	C-O	-6.03	1.16	1.23
1	A	87	VAL	CA-CB	6.00	1.60	1.53
1	B	335	GLY	C-O	5.88	1.27	1.24
1	A	319	VAL	CA-CB	5.71	1.61	1.54
1	B	378	ILE	CA-CB	5.50	1.60	1.54
1	A	605	ILE	CA-CB	5.44	1.59	1.53
1	B	362	ILE	CA-CB	5.21	1.62	1.54
1	B	727	VAL	CA-CB	5.06	1.60	1.54

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	660	GLN	N-CA-C	9.05	130.08	110.80
1	A	114	ASN	CB-CA-C	-7.20	101.52	111.95
1	A	354	ASN	N-CA-C	6.39	118.88	109.42
1	A	509	ASP	CA-C-N	-6.31	113.48	119.85
1	A	509	ASP	C-N-CA	-6.31	113.48	119.85
1	A	370	VAL	CB-CA-C	-5.99	100.35	111.79
1	B	328	GLY	N-CA-C	-5.83	100.45	112.34
1	B	68	ASN	N-CA-C	5.75	119.47	112.23
1	A	423	GLY	N-CA-C	-5.70	107.52	114.92
1	A	247	TRP	N-CA-C	-5.56	106.17	113.12
1	B	331	VAL	N-CA-C	-5.52	107.06	111.81
1	B	509	ASP	N-CA-C	5.48	115.90	109.60
1	B	335	GLY	N-CA-C	-5.46	105.54	111.21
1	A	468	ASN	CA-C-N	-5.24	114.52	119.76
1	A	468	ASN	C-N-CA	-5.24	114.52	119.76
1	B	318	ASN	CB-CA-C	-5.09	105.30	111.82
1	A	328	GLY	N-CA-C	-5.07	102.00	112.34
1	A	64	GLN	N-CA-C	5.04	117.43	111.33
1	A	328	GLY	CA-C-N	-5.01	115.40	120.31
1	A	328	GLY	C-N-CA	-5.01	115.40	120.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	659	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	5611	0	5302	44	0
1	B	5611	0	5300	48	0
2	C	39	0	34	0	0
2	D	39	0	34	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	12	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	12	0	10	0	0
5	A	609	0	0	5	0
5	B	584	0	0	3	0
All	All	12519	0	10690	92	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 4.

All (92) 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:448:ILE:HD11	1:B:489:MET:HG3	1.58	0.84
1:B:547:ASN:HD22	1:B:550:PHE:HE2	1.32	0.78
1:B:505:TRP:HB3	1:B:558:LEU:HD23	1.69	0.73
1:B:257:ASN:HD22	1:B:289:THR:HB	1.59	0.67
1:A:616:GLN:HG3	1:A:632:THR:HB	1.79	0.64
1:A:176:GLN:H	1:A:176:GLN:HE21	1.45	0.63
1:A:441:GLN:NE2	1:A:474:GLN:OE1	2.25	0.62
1:A:418:PHE:HE1	1:A:448:ILE:HD11	1.63	0.61
1:B:381:LYS:HG2	1:B:520:TRP:HZ2	1.67	0.60
1:B:616:GLN:HG3	1:B:632:THR:HB	1.83	0.60
1:A:547:ASN:HD22	1:A:550:PHE:HE2	1.49	0.59
1:B:107:GLN:HG2	1:B:152:SER:HB3	1.83	0.59
1:B:229:VAL:CG1	1:B:232:LEU:HD22	2.33	0.58
1:B:229:VAL:HG11	1:B:232:LEU:HD22	1.87	0.57
1:A:194:GLN:NE2	5:A:1331:HOH:O	2.38	0.56
1:B:455:VAL:HG22	1:B:490:ILE:HB	1.88	0.54
1:A:235:ASN:HD22	1:A:257:ASN:HB2	1.73	0.54
1:A:623:ALA:HA	1:A:672:HIS:O	2.08	0.54
1:B:231:ASN:HA	1:B:253:ARG:O	2.07	0.53
1:A:1:LEU:N	5:A:777:HOH:O	2.41	0.53
1:A:487:THR:HG22	1:A:520:TRP:HB2	1.90	0.53
1:B:381:LYS:HG2	1:B:520:TRP:CZ2	2.44	0.53
1:B:95:LYS:HE2	5:B:1044:HOH:O	2.10	0.52
1:A:246:ASN:OD1	1:A:246:ASN:C	2.54	0.51
1:B:448:ILE:CD1	1:B:489:MET:HG3	2.35	0.51
1:B:623:ALA:HA	1:B:672:HIS:O	2.11	0.51
1:A:427:ILE:HB	1:A:448:ILE:HD13	1.93	0.51
1:A:455:VAL:HG22	1:A:490:ILE:HB	1.94	0.49
1:A:381:LYS:HD3	1:A:520:TRP:HZ2	1.77	0.49
1:B:186:GLN:HE21	1:B:186:GLN:HA	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:VAL:HA	1:A:586:ILE:HG22	1.94	0.49
1:A:613:ILE:O	1:A:663:ALA:HB2	2.11	0.49
1:B:196:GLN:HA	1:B:218:ILE:O	2.13	0.48
1:B:702:ILE:HG23	1:B:703:VAL:HG23	1.95	0.48
1:A:381:LYS:HD3	1:A:520:TRP:CZ2	2.49	0.47
1:B:67:ILE:O	1:B:72:ARG:HD3	2.15	0.47
1:B:554:LEU:HD23	1:B:557:HIS:HB2	1.98	0.46
1:A:702:ILE:HG23	1:A:703:VAL:HG23	1.97	0.46
1:A:270:THR:HG23	1:A:306:GLN:HG3	1.98	0.46
1:B:613:ILE:CD1	1:B:646:PRO:HD2	2.45	0.46
1:A:418:PHE:CE1	1:A:448:ILE:HD11	2.49	0.46
1:A:186:GLN:HA	1:A:186:GLN:HE21	1.81	0.46
1:A:573:LEU:HD13	1:A:573:LEU:C	2.41	0.45
1:B:202:ASN:HB2	5:B:1133:HOH:O	2.15	0.45
1:B:317:THR:O	1:B:318:ASN:HB2	2.13	0.45
1:A:475:VAL:HG13	1:A:484:VAL:HG11	1.97	0.45
1:A:660:GLN:O	1:A:660:GLN:HG3	2.16	0.45
1:B:246:ASN:OD1	1:B:246:ASN:C	2.59	0.45
1:A:12:GLY:HA3	1:A:388:GLY:O	2.17	0.45
1:A:495:GLY:HA2	1:A:496:PRO:C	2.40	0.45
1:B:285:ASP:HB3	1:B:377:ARG:HD2	1.98	0.44
1:B:235:ASN:HD22	1:B:257:ASN:HB2	1.83	0.44
1:B:258:ASN:HA	1:B:290:ASN:O	2.18	0.44
1:A:483:VAL:HG22	1:A:516:ALA:HB1	1.99	0.44
1:B:536:VAL:HA	1:B:586:ILE:HG22	2.00	0.44
1:A:383:HIS:ND1	1:A:521:ASP:OD2	2.50	0.44
1:A:505:TRP:HB3	1:A:558:LEU:HD23	2.00	0.44
1:B:168:THR:HA	1:B:196:GLN:O	2.17	0.44
1:A:7:SER:HB2	1:A:8:PRO:CD	2.48	0.43
1:A:196:GLN:HA	1:A:218:ILE:O	2.18	0.43
1:B:160:LEU:HD12	1:B:187:MET:HG2	2.00	0.43
1:B:411:THR:HG22	1:B:415:LYS:HE2	2.00	0.43
1:A:145:ASN:HD22	1:A:203:GLY:HA2	1.84	0.43
1:B:154:ARG:HA	1:B:181:ILE:O	2.18	0.43
1:B:483:VAL:HG22	1:B:516:ALA:HB1	2.01	0.43
1:B:91:PRO:HG3	1:B:114:ASN:HB3	2.01	0.42
1:B:498:ALA:HB2	1:B:550:PHE:CG	2.54	0.42
1:A:316:LEU:O	1:A:336:THR:HA	2.20	0.42
1:A:67:ILE:O	1:A:72:ARG:HD3	2.19	0.42
1:B:317:THR:O	1:B:318:ASN:CB	2.67	0.42
1:B:284:ILE:HB	1:B:677:PHE:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:573:LEU:C	1:B:573:LEU:HD13	2.45	0.42
1:B:230:ARG:HA	1:B:252:GLN:O	2.19	0.42
1:A:75:GLN:HB2	5:A:1314:HOH:O	2.18	0.41
1:B:660:GLN:C	1:B:662:ASP:H	2.27	0.41
1:A:292:GLN:NE2	1:A:320:PRO:HG2	2.35	0.41
4:A:757:LGC:C1	5:A:1186:HOH:O	2.68	0.41
1:A:361:ASN:HD22	1:B:511:SER:HB3	1.86	0.41
1:B:462:LYS:O	1:B:462:LYS:HG3	2.20	0.41
1:A:263:PHE:HB2	1:A:295:VAL:HG22	2.02	0.41
1:A:553:PHE:O	1:A:591:ASN:HB3	2.21	0.41
1:B:475:VAL:HG13	1:B:484:VAL:HG11	2.03	0.41
1:A:174:VAL:HB	1:A:178:THR:HG21	2.03	0.41
1:A:481:SER:HB2	5:A:1227:HOH:O	2.21	0.41
1:A:253:ARG:HA	1:A:285:ASP:O	2.21	0.40
1:B:152:SER:HA	1:B:179:SER:O	2.22	0.40
1:B:316:LEU:O	1:B:336:THR:HA	2.21	0.40
1:B:683:SER:O	1:B:720:ASP:HB2	2.21	0.40
1:A:168:THR:HA	1:A:196:GLN:O	2.21	0.40
1:B:367:ARG:HB3	1:B:372:LEU:HD21	2.03	0.40
1:A:471:PRO:HA	1:A:501:ILE:O	2.21	0.40
1:B:30:ILE:HG22	5:B:866:HOH:O	2.21	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	750/758 (99%)	710 (95%)	38 (5%)	2 (0%)	36	40
1	B	750/758 (99%)	708 (94%)	38 (5%)	4 (0%)	24	24
All	All	1500/1516 (99%)	1418 (94%)	76 (5%)	6 (0%)	30	31

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	660	GLN
1	A	155	ASN
1	B	155	ASN
1	A	702	ILE
1	B	452	VAL
1	B	702	ILE

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	590/594 (99%)	571 (97%)	19 (3%)	34	43
1	B	590/594 (99%)	571 (97%)	19 (3%)	34	43
All	All	1180/1188 (99%)	1142 (97%)	38 (3%)	34	43

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	26	GLN
1	A	30	ILE
1	A	75	GLN
1	A	96	VAL
1	A	114	ASN
1	A	147	ASN
1	A	166	SER
1	A	176	GLN
1	A	186	GLN
1	A	227	PHE
1	A	229	VAL
1	A	238	ASN
1	A	302	SER
1	A	314	ILE
1	A	315	GLN
1	A	370	VAL

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Mol	Chain	Res	Type
1	A	427	ILE
1	A	660	GLN
1	B	10	THR
1	B	26	GLN
1	B	30	ILE
1	B	114	ASN
1	B	147	ASN
1	B	166	SER
1	B	186	GLN
1	B	227	PHE
1	B	271	SER
1	B	289	THR
1	B	315	GLN
1	B	318	ASN
1	B	448	ILE
1	B	466	TYR
1	B	573	LEU
1	B	658	LEU
1	B	660	GLN
1	B	707	SER
1	B	715	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	GLN
1	A	114	ASN
1	A	176	GLN
1	A	186	GLN
1	A	194	GLN
1	A	214	ASN
1	A	235	ASN
1	A	236	ASN
1	A	257	ASN
1	A	292	GLN
1	A	346	ASN
1	A	359	GLN
1	A	361	ASN
1	A	366	ASN
1	A	412	GLN
1	A	506	ASN
1	A	508	HIS

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Mol	Chain	Res	Type
1	A	535	GLN
1	A	547	ASN
1	A	631	GLN
1	A	672	HIS
1	A	696	GLN
1	A	737	ASN
1	A	738	ASN
1	B	23	GLN
1	B	114	ASN
1	B	186	GLN
1	B	214	ASN
1	B	217	ASN
1	B	235	ASN
1	B	238	ASN
1	B	242	ASN
1	B	257	ASN
1	B	290	ASN
1	B	313	ASN
1	B	359	GLN
1	B	547	ASN
1	B	642	ASN
1	B	752	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 ⓘ

6 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	1,2	14,14,15	0.82	0	17,19,21	1.86	4 (23%)
2	NAG	C	2	2	14,14,15	0.59	0	17,19,21	1.73	5 (29%)
2	BMA	C	3	2	11,11,12	1.26	1 (9%)	15,15,17	1.45	1 (6%)
2	NAG	D	1	1,2	14,14,15	0.77	0	17,19,21	1.42	2 (11%)
2	NAG	D	2	2	14,14,15	0.79	0	17,19,21	1.49	2 (11%)
2	BMA	D	3	2	11,11,12	0.87	0	15,15,17	1.38	2 (13%)

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	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	BMA	C	3	2	-	1/2/19/22	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	BMA	D	3	2	-	2/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	BMA	C2-C3	2.48	1.56	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	5.14	119.07	112.19
2	D	2	NAG	C1-O5-C5	4.66	118.44	112.19
2	C	3	BMA	C1-C2-C3	3.98	115.44	109.64
2	C	2	NAG	C1-O5-C5	3.31	116.62	112.19
2	D	1	NAG	C2-N2-C7	3.13	127.09	122.90
2	C	1	NAG	C2-N2-C7	-2.73	119.24	122.90
2	C	1	NAG	O5-C1-C2	-2.72	107.09	111.29
2	C	2	NAG	C2-N2-C7	-2.69	119.30	122.90
2	C	2	NAG	C4-C3-C2	2.54	114.74	111.02
2	D	1	NAG	O5-C1-C2	-2.51	107.40	111.29
2	C	1	NAG	C4-C3-C2	2.48	114.66	111.02
2	C	2	NAG	C8-C7-N2	2.44	120.17	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NAG	C1-C2-N2	-2.25	106.89	110.43
2	C	2	NAG	C3-C4-C5	-2.17	106.30	110.23
2	D	3	BMA	O5-C5-C4	-2.12	105.67	110.83
2	D	3	BMA	O5-C5-C6	2.02	111.59	107.66

There are no chirality outliers.

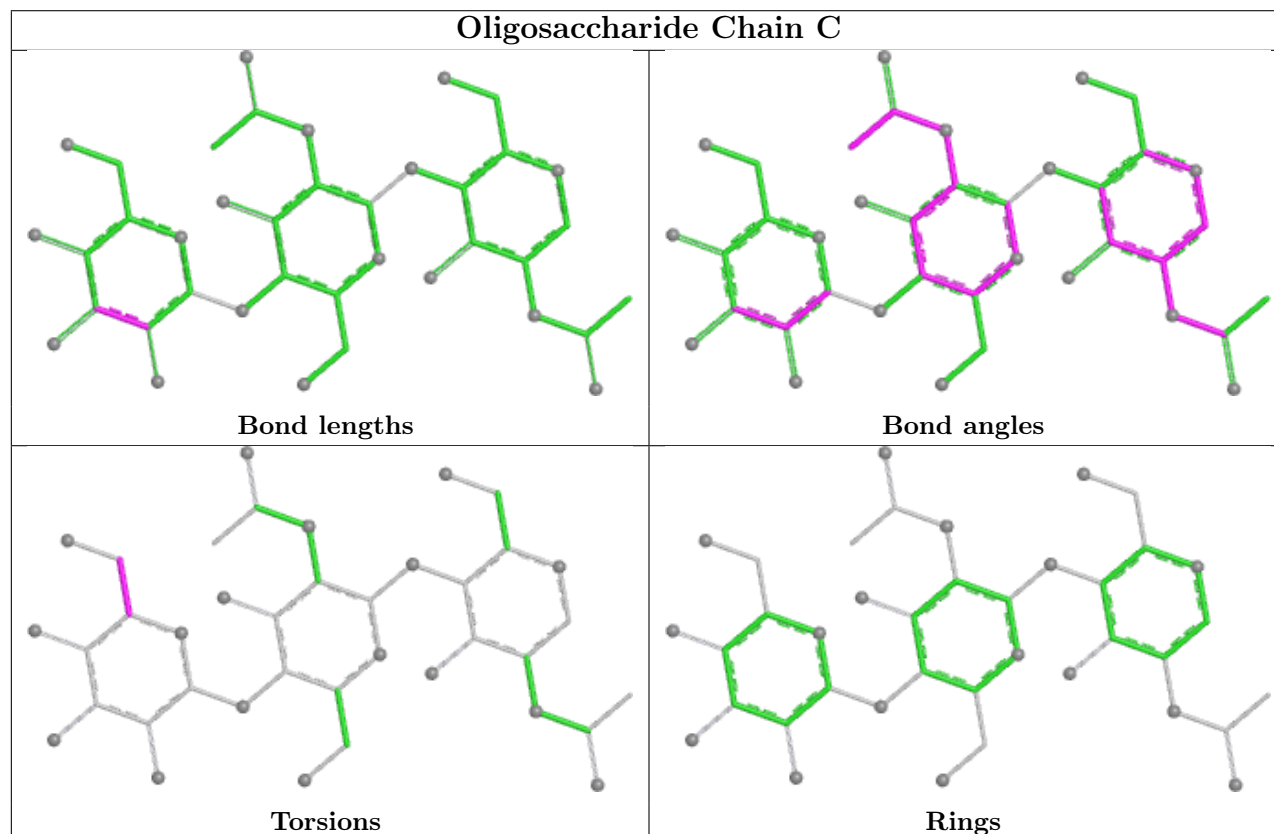
All (5) torsion outliers are listed below:

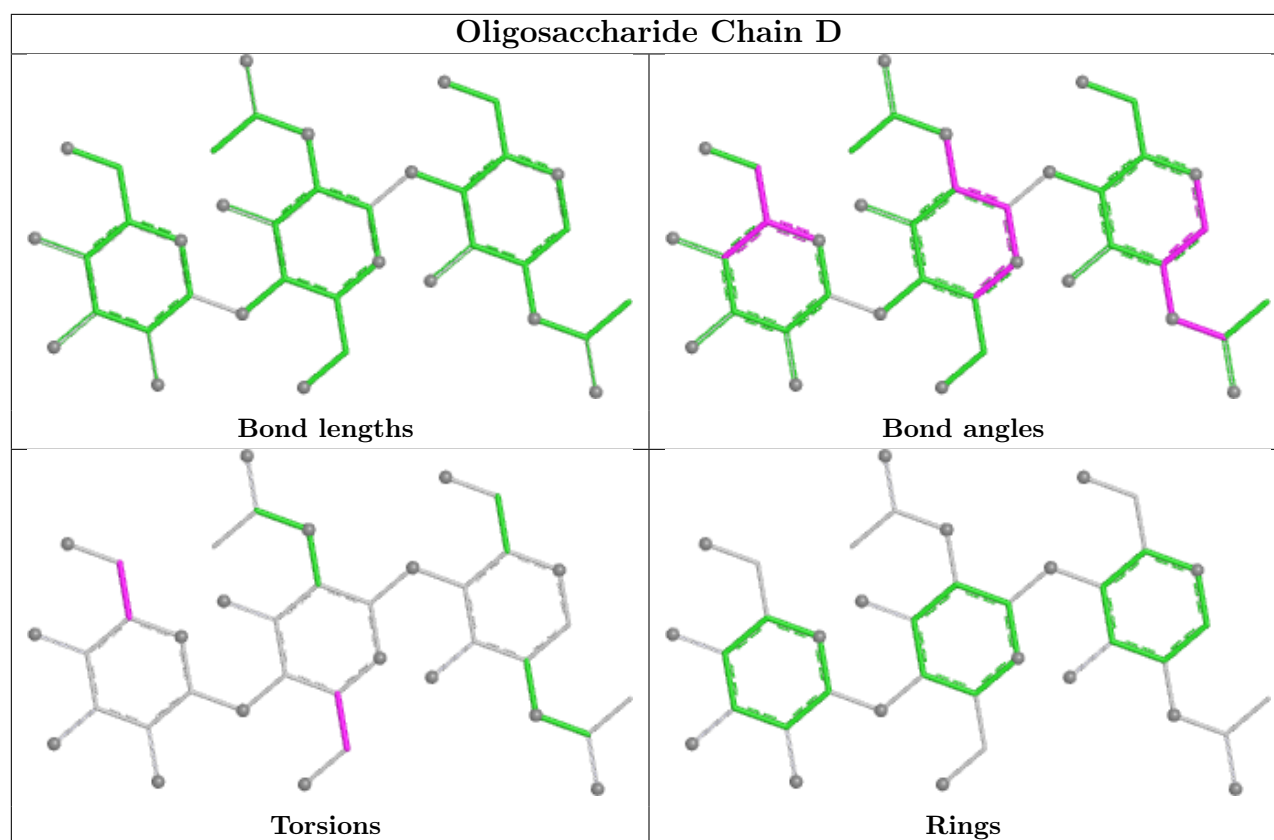
Mol	Chain	Res	Type	Atoms
2	D	3	BMA	O5-C5-C6-O6
2	D	3	BMA	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	C	3	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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	$\# Z > 2$
4	LGC	B	757	-	12,12,12	2.41	1 (8%)	15,17,17	1.32	2 (13%)
4	LGC	A	757	-	12,12,12	2.85	2 (16%)	15,17,17	1.37	2 (13%)

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	LGC	B	757	-	-	0/2/22/22	0/1/1/1
4	LGC	A	757	-	-	2/2/22/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	757	LGC	O5-C1	8.71	1.47	1.34
4	B	757	LGC	O5-C1	7.72	1.45	1.34
4	A	757	LGC	O5-C5	-3.93	1.41	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	757	LGC	O4-C4-C5	-2.72	102.63	109.32
4	A	757	LGC	O6-C6-C5	-2.60	102.49	111.33
4	A	757	LGC	O4-C4-C5	-2.40	103.40	109.32
4	B	757	LGC	O2-C2-C3	2.29	115.29	110.53

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	757	LGC	C4-C5-C6-O6
4	A	757	LGC	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	757	LGC	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	752/758 (99%)	-0.75	0	100 100	10, 18, 28, 41	0
1	B	752/758 (99%)	-0.66	1 (0%)	92 92	12, 20, 31, 48	0
All	All	1504/1516 (99%)	-0.71	1 (0%)	92 92	10, 19, 30, 48	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	5	CYS	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](#)

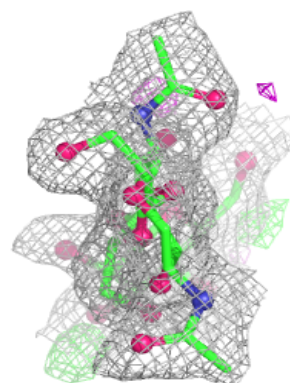
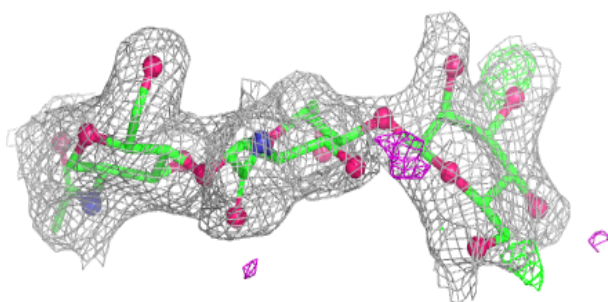
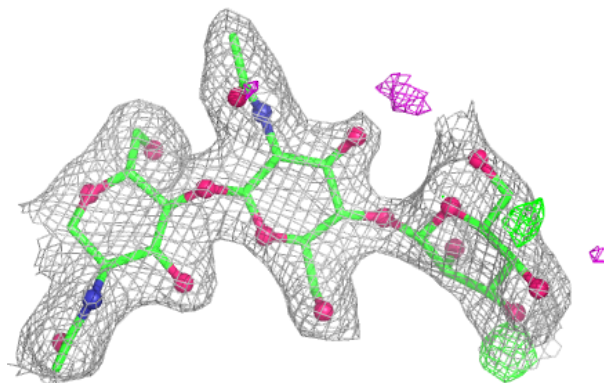
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BMA	C	3	11/12	0.67	0.16	60,62,63,63	0
2	BMA	D	3	11/12	0.73	0.14	65,69,71,72	0
2	NAG	C	2	14/15	0.86	0.11	41,44,50,56	0
2	NAG	D	2	14/15	0.89	0.10	31,43,52,60	0
2	NAG	D	1	14/15	0.95	0.05	20,24,29,36	0
2	NAG	C	1	14/15	0.96	0.06	21,28,30,36	0

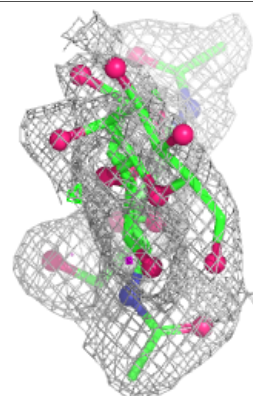
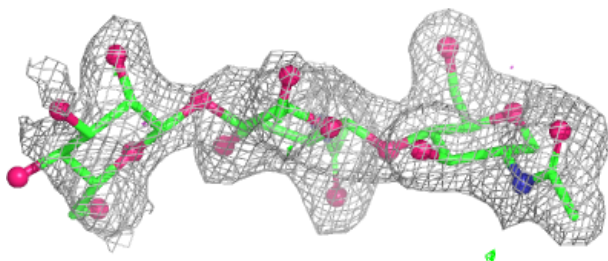
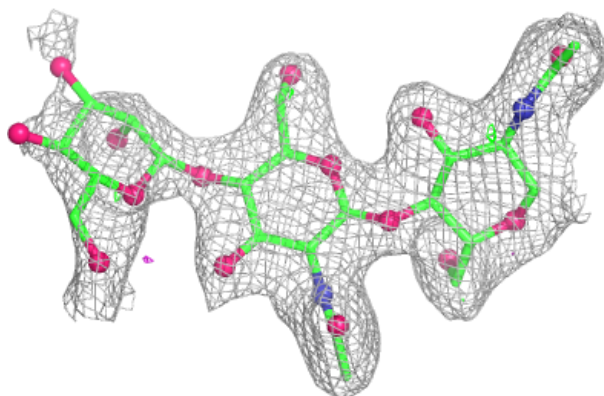
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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:**

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
4	LGC	A	757	12/12	0.95	0.06	16,24,28,28	0
4	LGC	B	757	12/12	0.95	0.06	21,26,30,31	0
3	ZN	A	756	1/1	0.97	0.04	38,38,38,38	0
3	ZN	B	756	1/1	0.99	0.03	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.