



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:38 PM UTC

PDB ID : 2O9R / pdb_00002o9r
Title : beta-glucosidase B complexed with thiocellobiose
Authors : Isorna, P.; Polaina, J.; Sanz-Aparicio, J.
Deposited on : 2006-12-14
Resolution : 2.30 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

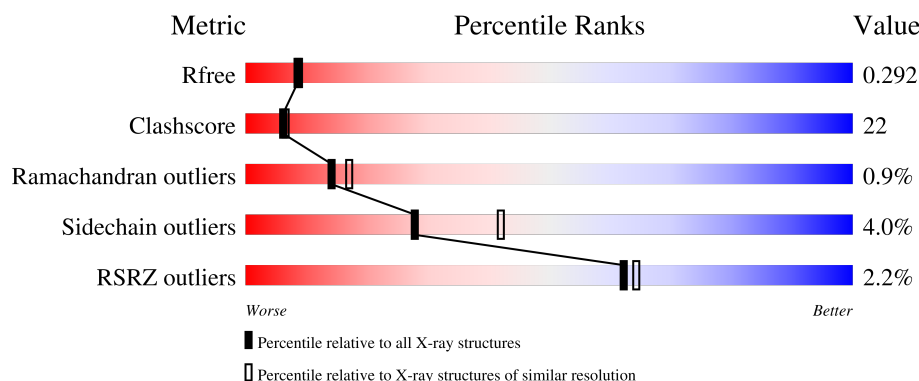
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

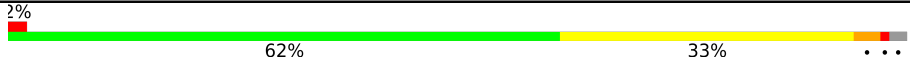
The reported resolution of this entry is 2.30 Å.

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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	452	 2% 62% 33% ...
2	B	2	 50% 50%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3778 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 Beta-glucosidase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	0	0
			3625	2327	612	666	20			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MET	-	expression tag	UNP P22505
A	-2	HIS	-	expression tag	UNP P22505
A	-1	HIS	-	expression tag	UNP P22505
A	0	HIS	-	expression tag	UNP P22505
A	1	HIS	-	expression tag	UNP P22505
A	2	HIS	-	expression tag	UNP P22505
A	3	HIS	-	expression tag	UNP P22505
A	376	GLN	HIS	engineered mutation	UNP P22505
A	377	ARG	GLY	engineered mutation	UNP P22505

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	S	0	0	0
			23	12	10	1			

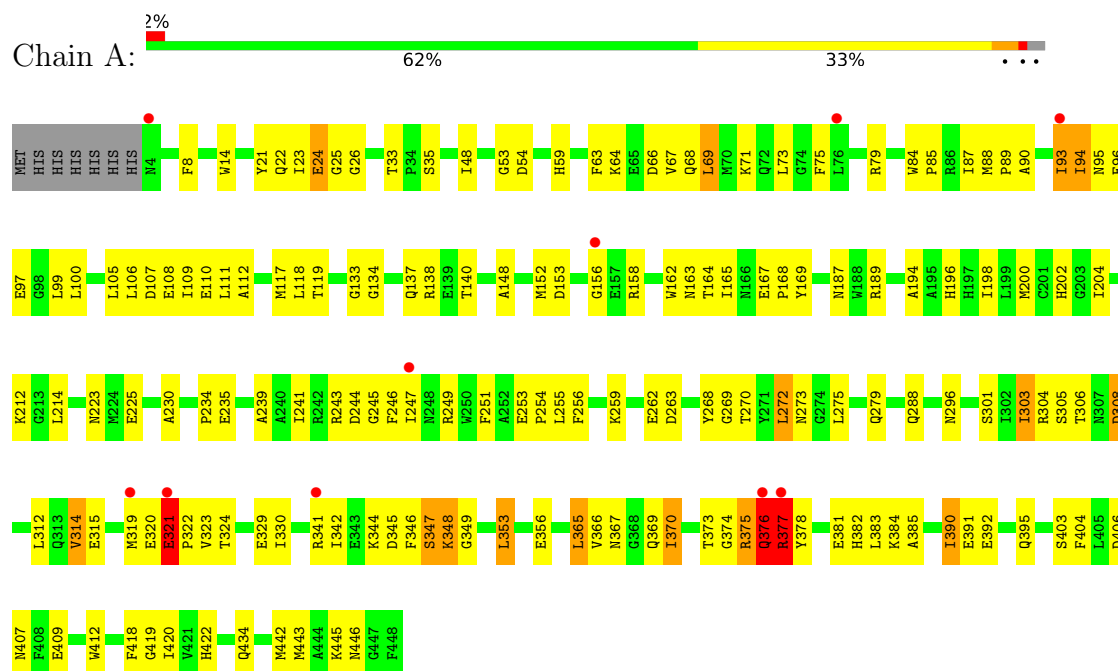
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	130	Total	O	0	0
			130	130		

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: Beta-glucosidase B



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.14Å 75.39Å 88.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.26 – 2.30 17.26 – 2.30	Depositor EDS
% Data completeness (in resolution range)	80.7 (17.26-2.30) 98.4 (17.26-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.30Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.201 , 0.273 0.223 , 0.292	Depositor DCC
R_{free} test set	1474 reflections (6.87%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.644	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3778	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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, SGC

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	0.47	2/3736 (0.1%)	1.00	20/5066 (0.4%)

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	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	375	ARG	C-N	-8.54	1.22	1.33
1	A	377	ARG	C-N	6.32	1.41	1.33

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ARG	O-C-N	11.84	138.18	122.43
1	A	375	ARG	CA-C-N	-10.33	99.48	121.80
1	A	375	ARG	C-N-CA	-10.33	99.48	121.80
1	A	376	GLN	OE1-CD-NE2	-9.79	112.81	122.60
1	A	24	GLU	N-CA-C	9.39	121.11	111.07
1	A	420	ILE	N-CA-C	-7.47	104.03	113.22
1	A	406	ASP	N-CA-C	-7.45	99.99	110.50
1	A	303	ILE	N-CA-C	6.89	119.36	108.89
1	A	376	GLN	CG-CD-NE2	6.67	126.41	116.40
1	A	390	ILE	N-CA-C	-6.61	95.59	109.34
1	A	404	PHE	N-CA-C	-6.56	104.21	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	347	SER	N-CA-C	6.48	119.16	111.71
1	A	321	GLU	CA-C-N	-6.21	112.78	119.98
1	A	321	GLU	C-N-CA	-6.21	112.78	119.98
1	A	384	LYS	N-CA-C	-6.17	104.66	111.82
1	A	63	PHE	N-CA-C	6.09	117.92	111.28
1	A	445	LYS	N-CA-C	-5.17	106.81	113.02
1	A	377	ARG	CA-C-N	-5.13	112.95	122.60
1	A	377	ARG	C-N-CA	-5.13	112.95	122.60
1	A	308	ASP	N-CA-C	5.09	117.73	109.79

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	375	ARG	Mainchain
1	A	377	ARG	Mainchain

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	3625	0	3423	159	0
2	B	23	0	21	1	0
3	A	130	0	0	5	0
All	All	3778	0	3444	159	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 22.

All (159) 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:321:GLU:HB3	1:A:322:PRO:CD	1.92	0.98
1:A:90:ALA:O	1:A:93:ILE:HD12	1.70	0.90
1:A:324:THR:HG23	1:A:330:ILE:HD11	1.55	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:GLY:HA2	1:A:162:TRP:HZ2	1.38	0.88
1:A:153:ASP:OD2	1:A:212:LYS:HE3	1.78	0.84
1:A:93:ILE:HD13	1:A:94:ILE:H	1.47	0.79
1:A:14:TRP:HB3	1:A:443:MET:HE3	1.62	0.78
1:A:14:TRP:CB	1:A:443:MET:HE3	2.13	0.78
1:A:253:GLU:HB3	1:A:254:PRO:HD3	1.66	0.77
1:A:374:GLY:C	1:A:376:GLN:H	1.89	0.76
1:A:48:ILE:HD12	1:A:412:TRP:HA	1.68	0.75
1:A:321:GLU:HB3	1:A:322:PRO:HD3	1.66	0.74
1:A:93:ILE:CD1	1:A:94:ILE:H	2.03	0.71
1:A:303:ILE:HG23	1:A:314:VAL:HG13	1.71	0.71
1:A:374:GLY:C	1:A:376:GLN:N	2.45	0.71
1:A:323:VAL:HG12	1:A:329:GLU:HA	1.75	0.69
1:A:165:ILE:HG21	1:A:198:ILE:HD11	1.75	0.69
1:A:341:ARG:HG2	3:A:607:HOH:O	1.94	0.67
1:A:390:ILE:O	1:A:391:GLU:HB3	1.93	0.67
1:A:245:GLY:HA2	1:A:249:ARG:HB2	1.76	0.66
1:A:370:ILE:N	1:A:370:ILE:HD12	2.12	0.65
1:A:75:PHE:HA	1:A:443:MET:HE1	1.78	0.65
1:A:165:ILE:CG2	1:A:198:ILE:HD11	2.26	0.65
1:A:196:HIS:CD2	1:A:200:MET:HE2	2.31	0.65
1:A:249:ARG:HH22	1:A:263:ASP:CG	2.05	0.65
1:A:93:ILE:CD1	1:A:93:ILE:H	2.10	0.64
1:A:164:THR:OG1	1:A:202:HIS:HD2	1.80	0.64
1:A:168:PRO:HA	1:A:198:ILE:HD13	1.79	0.63
1:A:243:ARG:NH1	1:A:247:ILE:HD11	2.13	0.63
1:A:249:ARG:NH2	1:A:263:ASP:OD1	2.31	0.63
1:A:239:ALA:HB1	1:A:314:VAL:HG22	1.79	0.63
1:A:194:ALA:O	1:A:198:ILE:HG22	1.99	0.62
1:A:246:PHE:HB2	1:A:312:LEU:HD11	1.81	0.62
1:A:79:ARG:HH21	1:A:163:ASN:ND2	1.97	0.61
1:A:71:LYS:HD2	1:A:112:ALA:HB1	1.82	0.61
1:A:106:LEU:O	1:A:110:GLU:HG3	2.01	0.61
1:A:22:GLN:O	1:A:407:ASN:HB2	2.00	0.61
1:A:342:ILE:O	1:A:345:ASP:O	2.19	0.61
1:A:366:VAL:O	1:A:369:GLN:HG2	2.01	0.61
1:A:189:ARG:HH11	1:A:189:ARG:HG3	1.66	0.60
1:A:67:VAL:HG21	1:A:108:GLU:HG3	1.84	0.60
1:A:376:GLN:C	1:A:378:TYR:H	2.10	0.60
1:A:14:TRP:HB2	1:A:443:MET:HE3	1.86	0.58
1:A:168:PRO:HD2	3:A:606:HOH:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:ILE:C	1:A:392:GLU:H	2.11	0.58
1:A:107:ASP:O	1:A:111:LEU:HD13	2.04	0.57
1:A:196:HIS:O	1:A:200:MET:HG3	2.04	0.57
1:A:202:HIS:HE1	1:A:288:GLN:O	1.88	0.57
1:A:376:GLN:C	1:A:378:TYR:N	2.62	0.57
1:A:64:LYS:HE2	1:A:68:GLN:NE2	2.20	0.56
1:A:148:ALA:O	1:A:152:MET:HG3	2.06	0.56
1:A:243:ARG:CZ	1:A:247:ILE:HD11	2.36	0.56
1:A:89:PRO:HD2	1:A:93:ILE:HD11	1.87	0.56
1:A:93:ILE:O	1:A:94:ILE:HB	2.05	0.56
1:A:321:GLU:CB	1:A:322:PRO:CD	2.78	0.56
1:A:324:THR:CG2	1:A:330:ILE:HD11	2.33	0.56
1:A:235:GLU:OE2	3:A:504:HOH:O	2.18	0.55
1:A:138:ARG:CZ	1:A:279:GLN:HG3	2.37	0.54
1:A:320:GLU:O	1:A:321:GLU:O	2.25	0.54
1:A:138:ARG:NH1	1:A:200:MET:HE1	2.23	0.54
1:A:347:SER:O	1:A:348:LYS:C	2.50	0.54
1:A:93:ILE:HD13	1:A:94:ILE:N	2.18	0.53
1:A:69:LEU:HD22	1:A:69:LEU:O	2.08	0.53
1:A:93:ILE:CG1	1:A:94:ILE:N	2.72	0.53
1:A:296:ASN:CG	1:A:356:GLU:HB2	2.34	0.53
1:A:167:GLU:CD	1:A:223:ASN:HB3	2.33	0.53
1:A:64:LYS:HA	1:A:108:GLU:HG2	1.90	0.53
1:A:268:TYR:HB3	1:A:272:LEU:HD13	1.91	0.53
1:A:303:ILE:HD12	1:A:314:VAL:CG1	2.39	0.52
1:A:93:ILE:CG1	1:A:94:ILE:H	2.23	0.52
1:A:93:ILE:HD13	1:A:93:ILE:H	1.74	0.51
1:A:304:ARG:HD2	1:A:315:GLU:OE2	2.10	0.51
1:A:390:ILE:HG22	1:A:391:GLU:H	1.75	0.51
1:A:69:LEU:O	1:A:73:LEU:HB2	2.11	0.51
1:A:378:TYR:CD1	1:A:378:TYR:C	2.89	0.51
1:A:84:TRP:HB3	1:A:85:PRO:HD3	1.92	0.51
1:A:321:GLU:O	1:A:323:VAL:HG13	2.10	0.51
1:A:200:MET:O	1:A:204:ILE:HG12	2.11	0.50
1:A:79:ARG:NH2	1:A:163:ASN:ND2	2.58	0.50
1:A:321:GLU:CD	1:A:322:PRO:HD3	2.36	0.50
1:A:321:GLU:OE2	1:A:322:PRO:HD3	2.11	0.50
1:A:319:MET:HE3	1:A:329:GLU:OE2	2.12	0.50
1:A:244:ASP:OD1	1:A:249:ARG:HD3	2.12	0.49
1:A:409:GLU:OE2	2:B:2:BGC:H6C1	2.12	0.49
1:A:64:LYS:HE2	1:A:68:GLN:HE22	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:VAL:HG23	1:A:68:GLN:N	2.28	0.49
1:A:382:HIS:O	1:A:385:ALA:HB3	2.13	0.49
1:A:230:ALA:N	1:A:303:ILE:O	2.43	0.48
1:A:321:GLU:HB3	1:A:322:PRO:HD2	1.85	0.48
1:A:249:ARG:O	1:A:253:GLU:HB2	2.13	0.48
1:A:243:ARG:NH2	3:A:605:HOH:O	2.46	0.48
1:A:87:ILE:HD12	1:A:99:LEU:HD23	1.94	0.48
1:A:93:ILE:HG12	1:A:94:ILE:N	2.28	0.48
1:A:133:GLY:HA3	1:A:137:GLN:HG2	1.96	0.47
1:A:442:MET:HE3	1:A:443:MET:HG3	1.96	0.47
1:A:377:ARG:O	1:A:381:GLU:HG2	2.15	0.47
1:A:259:LYS:HD2	1:A:262:GLU:OE2	2.14	0.47
1:A:403:SER:O	1:A:419:GLY:HA2	2.14	0.47
1:A:88:MET:HA	1:A:93:ILE:HD11	1.97	0.47
1:A:349:GLY:O	1:A:395:GLN:HG3	2.15	0.47
1:A:169:TYR:CE1	1:A:247:ILE:HG12	2.50	0.46
1:A:256:PHE:CD1	1:A:346:PHE:HB3	2.50	0.46
1:A:156:GLY:HA2	1:A:162:TRP:CZ2	2.31	0.46
1:A:93:ILE:HD13	1:A:93:ILE:N	2.30	0.46
1:A:344:LYS:HE3	1:A:345:ASP:OD1	2.15	0.46
1:A:25:GLY:O	1:A:26:GLY:C	2.59	0.46
1:A:369:GLN:C	1:A:370:ILE:HD12	2.41	0.46
1:A:306:THR:CG2	1:A:315:GLU:HB2	2.45	0.46
1:A:84:TRP:N	1:A:85:PRO:CD	2.80	0.45
1:A:255:LEU:HD23	1:A:256:PHE:CE1	2.50	0.45
1:A:8:PHE:O	1:A:446:ASN:ND2	2.49	0.45
1:A:96:GLU:O	1:A:100:LEU:HG	2.17	0.45
1:A:442:MET:C	1:A:442:MET:SD	3.00	0.45
1:A:95:ASN:ND2	1:A:97:GLU:H	2.14	0.45
1:A:134:GLY:O	1:A:140:THR:OG1	2.34	0.45
1:A:169:TYR:HA	1:A:251:PHE:CZ	2.51	0.45
1:A:321:GLU:CB	1:A:322:PRO:HD3	2.40	0.45
1:A:370:ILE:O	1:A:434:GLN:HG2	2.17	0.45
1:A:79:ARG:NH1	1:A:356:GLU:OE1	2.49	0.45
1:A:89:PRO:HG2	1:A:93:ILE:HG13	1.99	0.45
1:A:105:LEU:O	1:A:109:ILE:HG13	2.16	0.45
1:A:366:VAL:O	1:A:367:ASN:C	2.60	0.44
1:A:119:THR:HG23	1:A:163:ASN:HD22	1.82	0.44
1:A:225:GLU:OE2	1:A:301:SER:HB2	2.16	0.44
1:A:23:ILE:HD12	1:A:59:HIS:CG	2.53	0.44
1:A:119:THR:HA	1:A:163:ASN:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:TYR:CB	1:A:272:LEU:HD13	2.47	0.43
1:A:234:PRO:HD2	3:A:504:HOH:O	2.18	0.43
1:A:93:ILE:O	1:A:94:ILE:CB	2.67	0.43
1:A:308:ASP:C	1:A:308:ASP:OD2	2.61	0.43
1:A:403:SER:HB3	1:A:418:PHE:O	2.18	0.43
1:A:33:THR:OG1	1:A:89:PRO:HA	2.18	0.43
1:A:35:SER:HB3	1:A:85:PRO:HG2	2.00	0.43
1:A:370:ILE:N	1:A:370:ILE:CD1	2.81	0.43
1:A:306:THR:HG23	1:A:315:GLU:HB2	2.01	0.43
1:A:390:ILE:O	1:A:392:GLU:N	2.50	0.43
1:A:117:MET:HE1	1:A:353:LEU:HD23	2.01	0.42
1:A:373:THR:O	1:A:376:GLN:HB3	2.19	0.42
1:A:79:ARG:NH2	1:A:163:ASN:HD21	2.16	0.42
1:A:23:ILE:HG13	1:A:24:GLU:N	2.35	0.42
1:A:71:LYS:HD2	1:A:112:ALA:CB	2.47	0.42
1:A:21:TYR:CZ	1:A:53:GLY:HA3	2.54	0.42
1:A:187:ASN:OD1	1:A:189:ARG:HB3	2.19	0.42
1:A:365:LEU:HD12	1:A:422:HIS:CE1	2.55	0.42
1:A:329:GLU:C	1:A:330:ILE:HD13	2.45	0.42
1:A:330:ILE:HD12	1:A:378:TYR:CE2	2.56	0.41
1:A:189:ARG:NH1	1:A:273:ASN:HB2	2.36	0.41
1:A:243:ARG:O	1:A:243:ARG:HD3	2.20	0.41
1:A:118:LEU:C	1:A:118:LEU:HD23	2.45	0.41
1:A:165:ILE:HG21	1:A:198:ILE:CD1	2.46	0.41
1:A:303:ILE:HD12	1:A:314:VAL:HG13	2.02	0.41
1:A:383:LEU:HD23	1:A:383:LEU:HA	1.96	0.41
1:A:89:PRO:CD	1:A:93:ILE:HD11	2.52	0.40
1:A:152:MET:HB3	1:A:214:LEU:HD11	2.03	0.40
1:A:390:ILE:C	1:A:392:GLU:N	2.73	0.40
1:A:59:HIS:HD2	1:A:66:ASP:OD2	2.03	0.40
1:A:246:PHE:CB	1:A:312:LEU:HD11	2.49	0.40
1:A:390:ILE:O	1:A:391:GLU:CB	2.62	0.40
1:A:71:LYS:HB2	1:A:112:ALA:HB1	2.04	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	443/452 (98%)	410 (93%)	29 (6%)	4 (1%)	14 17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	321	GLU
1	A	94	ILE
1	A	269	GLY
1	A	348	LYS

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	377/384 (98%)	362 (96%)	15 (4%)	28 42

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	69	LEU
1	A	93	ILE
1	A	158	ARG
1	A	241	ILE
1	A	270	THR
1	A	272	LEU

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Mol	Chain	Res	Type
1	A	275	LEU
1	A	305	SER
1	A	314	VAL
1	A	321	GLU
1	A	353	LEU
1	A	365	LEU
1	A	370	ILE
1	A	376	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	42	GLN
1	A	59	HIS
1	A	68	GLN
1	A	77	HIS
1	A	95	ASN
1	A	127	GLN
1	A	163	ASN
1	A	166	ASN
1	A	202	HIS
1	A	207	ASN
1	A	209	HIS
1	A	226	HIS
1	A	248	ASN
1	A	257	ASN
1	A	273	ASN
1	A	287	GLN
1	A	288	GLN
1	A	307	ASN
1	A	313	GLN
1	A	369	GLN
1	A	441	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 ⓘ

2 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	SGC	B	1	2	11,12,12	4.91	9 (81%)	14,17,17	1.63	2 (14%)
2	BGC	B	2	2	11,11,12	2.60	5 (45%)	15,15,17	2.18	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	SGC	B	1	2	-	2/2/22/22	0/1/1/1
2	BGC	B	2	2	-	2/2/19/22	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	SGC	C5-C4	10.46	1.61	1.53
2	B	1	SGC	C3-C4	9.53	1.61	1.53
2	B	2	BGC	C2-C3	4.90	1.60	1.52
2	B	2	BGC	O5-C1	3.54	1.49	1.43
2	B	2	BGC	C1-C2	3.36	1.60	1.52
2	B	2	BGC	C4-C5	3.34	1.60	1.53
2	B	1	SGC	O5-C1	3.30	1.50	1.42
2	B	1	SGC	O5-C5	3.22	1.52	1.44
2	B	1	SGC	O1-C1	3.18	1.49	1.39
2	B	1	SGC	C3-C2	3.05	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	SGC	O3-C3	2.88	1.50	1.43
2	B	1	SGC	C6-C5	2.72	1.60	1.51
2	B	1	SGC	C1-C2	2.45	1.58	1.52
2	B	2	BGC	C4-C3	2.44	1.58	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	BGC	C1-O5-C5	7.17	121.80	112.19
2	B	1	SGC	O3-C3-C2	-4.07	100.77	110.38
2	B	1	SGC	O3-C3-C4	3.02	115.34	109.17
2	B	2	BGC	C6-C5-C4	2.38	118.86	113.02
2	B	2	BGC	O5-C5-C4	-2.34	105.14	110.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

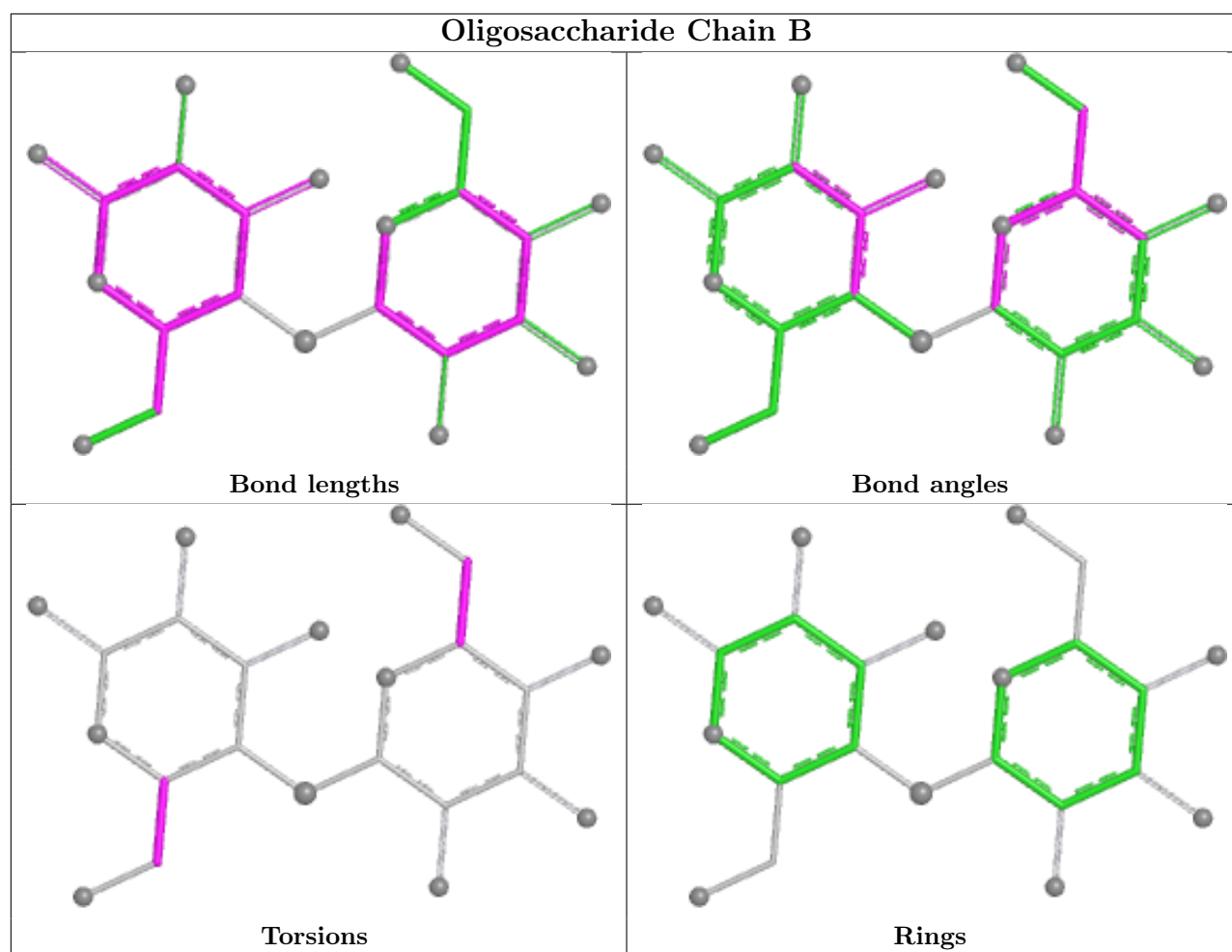
Mol	Chain	Res	Type	Atoms
2	B	1	SGC	O5-C5-C6-O6
2	B	1	SGC	C4-C5-C6-O6
2	B	2	BGC	O5-C5-C6-O6
2	B	2	BGC	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	BGC	1	0

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



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 [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	445/452 (98%)	0.09	10 (2%) 62 64	20, 33, 44, 56	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	376	GLN	3.7
1	A	319	MET	3.2
1	A	156	GLY	2.8
1	A	321	GLU	2.7
1	A	377	ARG	2.6
1	A	76	LEU	2.3
1	A	4	ASN	2.1
1	A	247	ILE	2.1
1	A	93	ILE	2.1
1	A	341	ARG	2.0

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

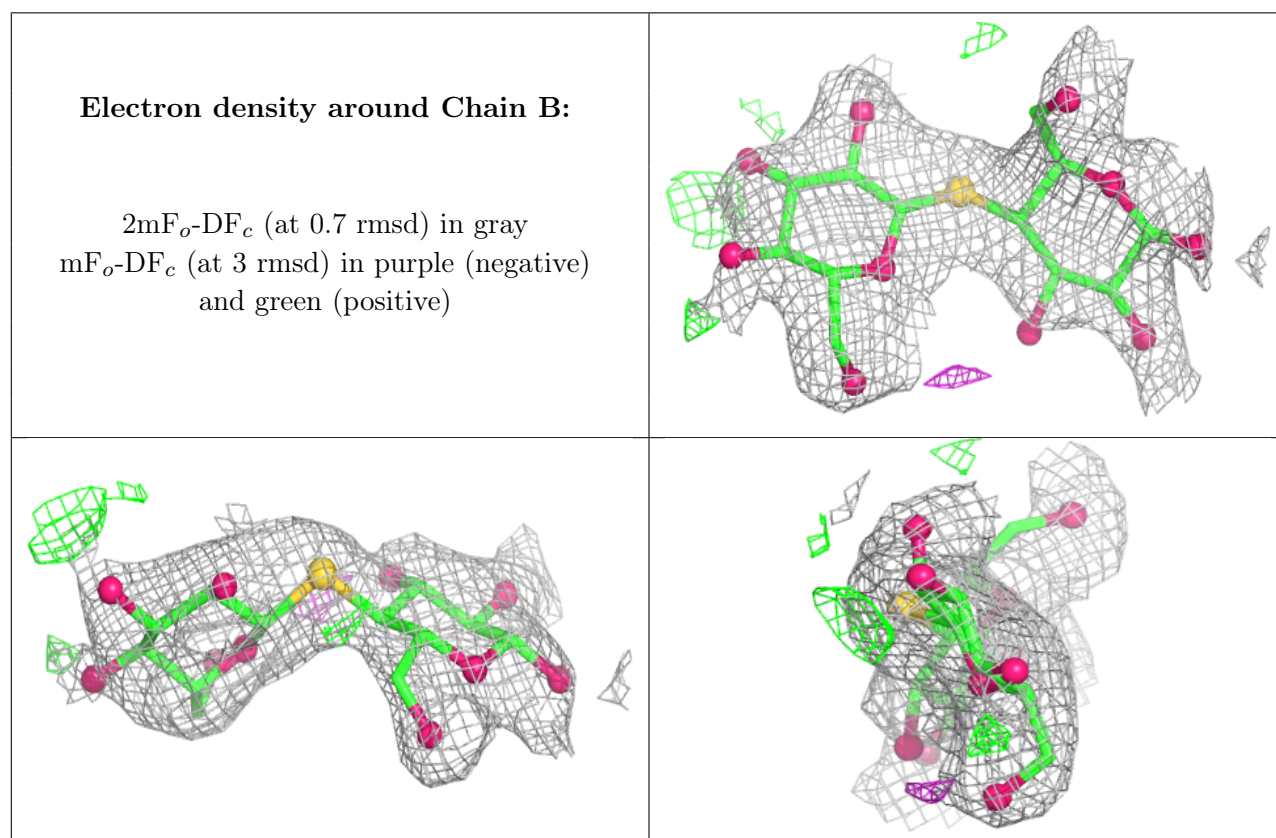
There are no non-standard protein/DNA/RNA residues in this entry.

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	SGC	B	1	12/12	0.77	0.13	55,56,57,58	0
2	BGC	B	2	11/12	0.84	0.12	48,51,52,53	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.



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.