



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 11:25 AM BST

PDB ID : 5NGY
Title : Crystal structure of Leuconostoc citreum NRRL B-1299 dextransucrase DSR-M
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Deposited on : 2017-03-20
Resolution : 3.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

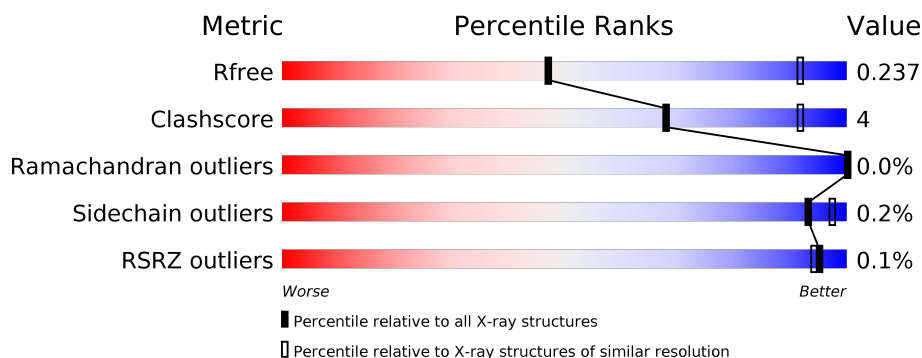
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1049 (3.88-3.52)
Clashscore	141614	1027 (3.86-3.54)
Ramachandran outliers	138981	1069 (3.88-3.52)
Sidechain outliers	138945	1065 (3.88-3.52)
RSRZ outliers	127900	1578 (3.90-3.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1290	
1	B	1290	
2	C	4	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	C	4	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 18858 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 DSR-M glucansucrase inactive mutant E715Q.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1265	Total	C	N	O	S	0	0	0
			9840	6167	1660	1994	19			
1	B	1144	Total	C	N	O	S	0	0	0
			8965	5618	1512	1816	19			

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	0	0	0
			45	24	21			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is PRASEODYMIUM ION (three-letter code: PR) (formula: Pr).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Pr	0	0
			3	3		
4	A	3	Total	Pr	0	0
			3	3		

- Molecule 2: alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose

Chain C: 


GLC1
GLC2
GLC3
GLC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.87Å 128.81Å 234.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.70 48.25 – 3.70	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-3.70) 98.2 (48.25-3.70)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 3.67Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.192 , 0.236 0.195 , 0.237	Depositor DCC
R_{free} test set	1693 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	62.3	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 4.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	18858	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.39	1/10056 (0.0%)	0.57	0/13672
1	B	0.38	0/9161	0.57	0/12448
All	All	0.38	1/19217 (0.0%)	0.57	0/26120

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	495	ALA	C-N	6.80	1.49	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9840	0	9208	81	0
1	B	8965	0	8456	60	0
2	C	45	0	39	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
All	All	18858	0	17703	139	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 (139) 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:990:PRO:HG2	1:A:993:LEU:HD21	1.58	0.86
1:A:297:PRO:O	1:A:298:ASP:HB3	1.82	0.79
1:A:990:PRO:HG2	1:A:993:LEU:CD2	2.13	0.78
1:B:329:PHE:CE1	1:B:334:ALA:HB2	2.19	0.77
1:A:261:LYS:O	1:A:276:ASN:O	2.05	0.74
1:A:351:ASP:OD1	1:A:352:LEU:N	2.23	0.70
1:B:351:ASP:OD1	1:B:355:ASN:HB2	1.94	0.68
1:A:1117:LEU:O	1:A:1123:VAL:HG11	1.97	0.65
1:A:297:PRO:O	1:A:298:ASP:CB	2.44	0.65
1:A:334:ALA:O	1:A:359:PHE:HE1	1.79	0.65
1:B:1117:LEU:O	1:B:1123:VAL:HG11	1.97	0.64
1:B:1402:TYR:CD2	1:B:1415:ILE:CD1	2.81	0.64
1:A:1182:ASP:HB3	1:A:1183:GLN:OE1	2.01	0.61
1:A:978:GLN:N	1:A:978:GLN:OE1	2.26	0.61
1:B:435:ASP:OD2	1:B:437:LYS:HE2	2.00	0.61
1:A:675:ARG:NH2	1:A:715:GLN:OE1	2.33	0.61
1:A:267:ASN:OD1	1:A:268:SER:N	2.33	0.61
1:B:1352:TYR:CD1	1:B:1372:VAL:HG21	2.36	0.60
1:B:1182:ASP:HB3	1:B:1183:GLN:OE1	2.01	0.60
1:A:1219:ILE:HG13	1:A:1219:ILE:O	2.02	0.60
1:A:644:GLU:HG3	1:A:1132:ARG:NH2	2.16	0.60
1:A:344:ILE:HG21	1:A:349:ALA:HB2	1.82	0.59
1:A:1373:ASN:OD1	1:A:1375:SER:HB3	2.03	0.59
1:B:1352:TYR:CE1	1:B:1372:VAL:CG2	2.86	0.58
1:B:1341:THR:HG22	1:B:1342:PHE:O	2.04	0.57
1:A:334:ALA:O	1:A:359:PHE:CE1	2.58	0.57
1:A:1341:THR:HG22	1:A:1342:PHE:O	2.04	0.56
1:A:1387:GLU:OE1	1:A:1389:ARG:NH2	2.40	0.55
1:A:1380:TYR:HB2	1:A:1403:PHE:CZ	2.42	0.55
1:A:243:THR:OG1	1:A:250:LEU:CB	2.56	0.54
1:A:565:ILE:CG1	1:A:1280:GLY:HA3	2.38	0.54
1:B:1387:GLU:OE1	1:B:1389:ARG:NH2	2.40	0.54
1:B:1402:TYR:CD2	1:B:1415:ILE:HD13	2.43	0.54
1:B:1416:SER:O	1:B:1417:ASP:HB3	2.08	0.54
1:A:397:PHE:CE2	1:A:1382:LEU:HD22	2.43	0.54
1:A:565:ILE:HG13	1:A:1280:GLY:HA3	1.89	0.54
1:B:1380:TYR:HB2	1:B:1403:PHE:CZ	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:GLY:O	1:B:333:THR:OG1	2.25	0.53
2:C:3:GLC:H62	2:C:4:GLC:O2	2.09	0.53
1:A:992:ASN:OD1	1:A:993:LEU:HD22	2.09	0.53
1:A:306:PHE:CE1	1:A:315:TYR:HB2	2.43	0.52
1:A:1324:ASP:HB3	1:A:1327:GLY:O	2.09	0.52
1:A:208:GLN:NE2	1:A:238:TYR:CE1	2.77	0.52
1:B:894:THR:HG22	1:B:1096:LYS:O	2.10	0.52
1:B:643:ALA:HA	1:B:1259:ILE:HD11	1.92	0.52
1:B:1417:ASP:OD1	1:B:1418:ASP:N	2.43	0.52
1:B:1324:ASP:HB3	1:B:1327:GLY:O	2.10	0.52
1:A:309:VAL:HG23	1:A:310:ASN:OD1	2.09	0.51
1:B:1123:VAL:HG13	1:B:1124:GLN:N	2.25	0.51
1:B:312:LYS:O	1:B:341:GLY:O	2.27	0.51
1:A:894:THR:HG22	1:A:1096:LYS:O	2.10	0.51
1:A:344:ILE:HG23	1:A:344:ILE:O	2.09	0.51
1:A:1123:VAL:HG13	1:A:1124:GLN:N	2.25	0.51
1:B:582:PHE:CE2	1:B:1219:ILE:CD1	2.94	0.51
1:B:976:ASN:HD21	1:B:981:VAL:CG2	2.24	0.51
1:A:643:ALA:HA	1:A:1259:ILE:HD11	1.92	0.50
1:B:975:THR:HG23	1:B:1022:LEU:O	2.12	0.50
1:A:990:PRO:CG	1:A:993:LEU:CD2	2.89	0.50
1:B:1410:VAL:HG13	1:B:1413:GLN:NE2	2.27	0.50
1:A:397:PHE:CE2	1:A:1382:LEU:CD2	2.95	0.50
1:A:308:GLN:HA	1:A:313:ASN:H	1.76	0.49
1:A:702:GLU:OE2	1:B:1371:THR:HG21	2.12	0.49
1:B:1219:ILE:HG12	1:B:1219:ILE:O	2.13	0.49
1:A:306:PHE:HB3	1:A:313:ASN:OD1	2.12	0.48
1:A:291:LYS:CB	1:A:322:TYR:CD2	2.95	0.48
1:A:320:THR:HB	1:A:322:TYR:CD1	2.49	0.48
1:A:382:GLN:HE22	1:A:1434:LYS:HB2	1.78	0.48
1:B:821:SER:HB3	1:B:1122:THR:HG21	1.96	0.48
1:A:346:SER:OG	1:A:361:ALA:N	2.47	0.48
1:A:382:GLN:NE2	1:A:1434:LYS:HB2	2.28	0.48
1:A:208:GLN:NE2	1:A:238:TYR:HE1	2.11	0.48
1:A:179:ARG:CB	1:A:753:THR:HG21	2.44	0.47
1:A:276:ASN:HD21	1:A:280:GLU:HB2	1.79	0.47
1:A:644:GLU:HG3	1:A:1132:ARG:HH22	1.79	0.47
1:B:1352:TYR:CE1	1:B:1372:VAL:HG21	2.49	0.47
1:B:346:SER:HA	1:B:359:PHE:O	2.15	0.47
1:A:382:GLN:OE1	1:A:382:GLN:N	2.47	0.47
1:B:1402:TYR:CD2	1:B:1415:ILE:HD11	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:655:PHE:CE1	1:B:659:THR:HG21	2.50	0.47
1:A:1352:TYR:CD1	1:A:1372:VAL:HG21	2.50	0.47
1:A:821:SER:HB3	1:A:1122:THR:HG21	1.96	0.47
1:B:296:GLY:O	1:B:298:ASP:N	2.47	0.47
1:A:1410:VAL:HG13	1:A:1413:GLN:NE2	2.29	0.47
1:A:655:PHE:CE1	1:A:659:THR:HG21	2.50	0.46
1:B:434:TYR:CE1	1:B:1313:MET:HB3	2.50	0.46
1:B:867:TYR:CE2	1:B:869:GLY:HA3	2.51	0.46
1:A:975:THR:HG23	1:A:1022:LEU:O	2.17	0.45
1:B:1120:VAL:HG22	1:B:1123:VAL:HG12	1.99	0.45
1:B:759:LYS:HA	1:B:908:VAL:HG21	1.98	0.45
1:A:285:LEU:HD23	1:A:293:GLN:O	2.17	0.44
1:A:1120:VAL:HG22	1:A:1123:VAL:HG12	1.98	0.44
1:A:1123:VAL:O	1:A:1148:LYS:HD2	2.18	0.44
1:A:1324:ASP:OD1	1:A:1325:ASP:N	2.50	0.44
1:A:1044:LYS:HD2	1:B:1378:LYS:HE2	2.00	0.44
1:A:265:VAL:HG22	1:A:275:PHE:HE1	1.83	0.44
1:A:348:ILE:N	1:A:348:ILE:HD12	2.33	0.44
1:B:1123:VAL:O	1:B:1148:LYS:HD2	2.18	0.44
1:A:867:TYR:CE2	1:A:869:GLY:HA3	2.52	0.43
1:A:759:LYS:HA	1:A:908:VAL:HG21	2.00	0.43
1:B:1324:ASP:OD1	1:B:1325:ASP:N	2.50	0.43
1:B:603:THR:HG21	1:B:625:GLY:HA3	2.00	0.43
1:B:947:PRO:HG3	1:B:1019:SER:OG	2.19	0.43
1:A:603:THR:HG21	1:A:625:GLY:HA3	2.00	0.43
1:B:320:THR:HB	1:B:322:TYR:CD1	2.53	0.43
1:A:310:ASN:OD1	1:A:312:LYS:HG2	2.18	0.43
1:A:819:LYS:HB3	1:A:1120:VAL:HG11	2.00	0.43
1:B:1219:ILE:CG1	1:B:1219:ILE:O	2.66	0.43
1:A:1072:VAL:HG23	1:A:1122:THR:O	2.19	0.43
1:A:797:ILE:HD11	1:A:828:PHE:CE1	2.54	0.43
1:B:1072:VAL:HG23	1:B:1122:THR:O	2.19	0.43
1:A:241:ASP:HB3	1:A:265:VAL:HB	2.01	0.43
1:B:303:LYS:NZ	1:B:303:LYS:HB2	2.34	0.42
1:A:1123:VAL:HG13	1:A:1124:GLN:H	1.84	0.42
1:A:460:ASN:HD22	1:A:529:SER:HB2	1.85	0.42
1:B:306:PHE:CE1	1:B:315:TYR:HB2	2.55	0.42
1:A:287:THR:HG22	1:A:292:THR:HG22	2.01	0.42
1:A:947:PRO:HG3	1:A:1019:SER:OG	2.20	0.42
1:B:645:GLN:OE1	1:B:1132:ARG:NH2	2.53	0.42
1:A:1380:TYR:HB2	1:A:1403:PHE:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:GLN:OE1	1:A:217:GLN:NE2	2.50	0.42
1:B:329:PHE:HB3	1:B:331:ASP:O	2.20	0.42
1:A:1049:SER:OG	1:A:1051:TYR:O	2.27	0.41
1:A:603:THR:HB	1:A:604:PRO:HD2	2.03	0.41
1:B:397:PHE:CE2	1:B:1382:LEU:HD22	2.55	0.41
1:B:603:THR:HB	1:B:604:PRO:HD2	2.02	0.41
1:B:787:ARG:NH1	1:B:870:ASP:OD1	2.44	0.41
1:B:797:ILE:HD11	1:B:828:PHE:CE1	2.55	0.41
1:A:848:ASN:HD21	1:A:1019:SER:HB3	1.86	0.41
1:B:1021:TYR:CE2	1:B:1023:ALA:HB2	2.55	0.41
1:B:380:ASP:HB3	1:B:386:ARG:HD2	2.02	0.41
1:B:848:ASN:HD21	1:B:1019:SER:HB3	1.85	0.41
1:B:648:TRP:CE3	1:B:1132:ARG:NH2	2.89	0.41
1:B:1380:TYR:HB2	1:B:1403:PHE:CE2	2.56	0.40
1:A:702:GLU:CD	1:B:1371:THR:HG21	2.42	0.40
1:B:1373:ASN:HA	1:B:1373:ASN:HD22	1.72	0.40
1:A:182:LEU:HD23	1:A:182:LEU:C	2.41	0.40
1:A:235:ASN:HB2	2:C:1:GLC:H2	2.04	0.40
1:A:291:LYS:CB	1:A:322:TYR:CE2	3.04	0.40
1:B:397:PHE:CE2	1:B:1382:LEU:CD2	3.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1263/1290 (98%)	1197 (95%)	65 (5%)	1 (0%)	51	83
1	B	1140/1290 (88%)	1085 (95%)	55 (5%)	0	100	100
All	All	2403/2580 (93%)	2282 (95%)	120 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	496	LEU

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	1043/1095 (95%)	1040 (100%)	3 (0%)	92	96
1	B	964/1095 (88%)	962 (100%)	2 (0%)	93	97
All	All	2007/2190 (92%)	2002 (100%)	5 (0%)	93	97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	THR
1	A	397	PHE
1	A	807	ASP
1	B	397	PHE
1	B	807	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	342	ASN
1	A	460	ASN
1	A	617	GLN
1	A	953	ASN
1	A	1145	ASN
1	A	1186	ASN
1	A	1212	ASN
1	A	1413	GLN
1	B	293	GLN
1	B	355	ASN
1	B	617	GLN
1	B	952	ASN
1	B	953	ASN

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Mol	Chain	Res	Type
1	B	976	ASN
1	B	1145	ASN
1	B	1186	ASN
1	B	1212	ASN
1	B	1373	ASN
1	B	1413	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 ⓘ

4 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	GLC	C	1	2	12,12,12	0.93	1 (8%)	17,17,17	1.99	4 (23%)
2	GLC	C	2	2	11,11,12	0.69	0	15,15,17	1.37	2 (13%)
2	GLC	C	3	2	11,11,12	0.59	0	15,15,17	1.06	1 (6%)
2	GLC	C	4	2	11,11,12	0.76	0	15,15,17	1.53	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	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	3	2	-	2/2/19/22	0/1/1/1
2	GLC	C	4	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	1	GLC	O1-C1	2.31	1.47	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	O1-C1-O5	5.88	128.03	110.38
2	C	4	GLC	C1-O5-C5	3.72	117.24	112.19
2	C	2	GLC	O5-C1-C2	-3.65	105.14	110.77
2	C	1	GLC	O5-C1-C2	2.98	115.61	110.28
2	C	3	GLC	C1-O5-C5	2.85	116.05	112.19
2	C	1	GLC	O2-C2-C1	2.78	115.60	109.16
2	C	2	GLC	O4-C4-C3	-2.59	104.36	110.35
2	C	1	GLC	C1-C2-C3	-2.29	105.56	110.31
2	C	4	GLC	O5-C1-C2	-2.18	107.41	110.77
2	C	4	GLC	O2-C2-C1	2.11	113.47	109.15

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	4	GLC	C4-C5-C6-O6
2	C	3	GLC	O5-C5-C6-O6
2	C	4	GLC	O5-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6
2	C	2	GLC	O5-C5-C6-O6
2	C	3	GLC	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 2 short contacts:

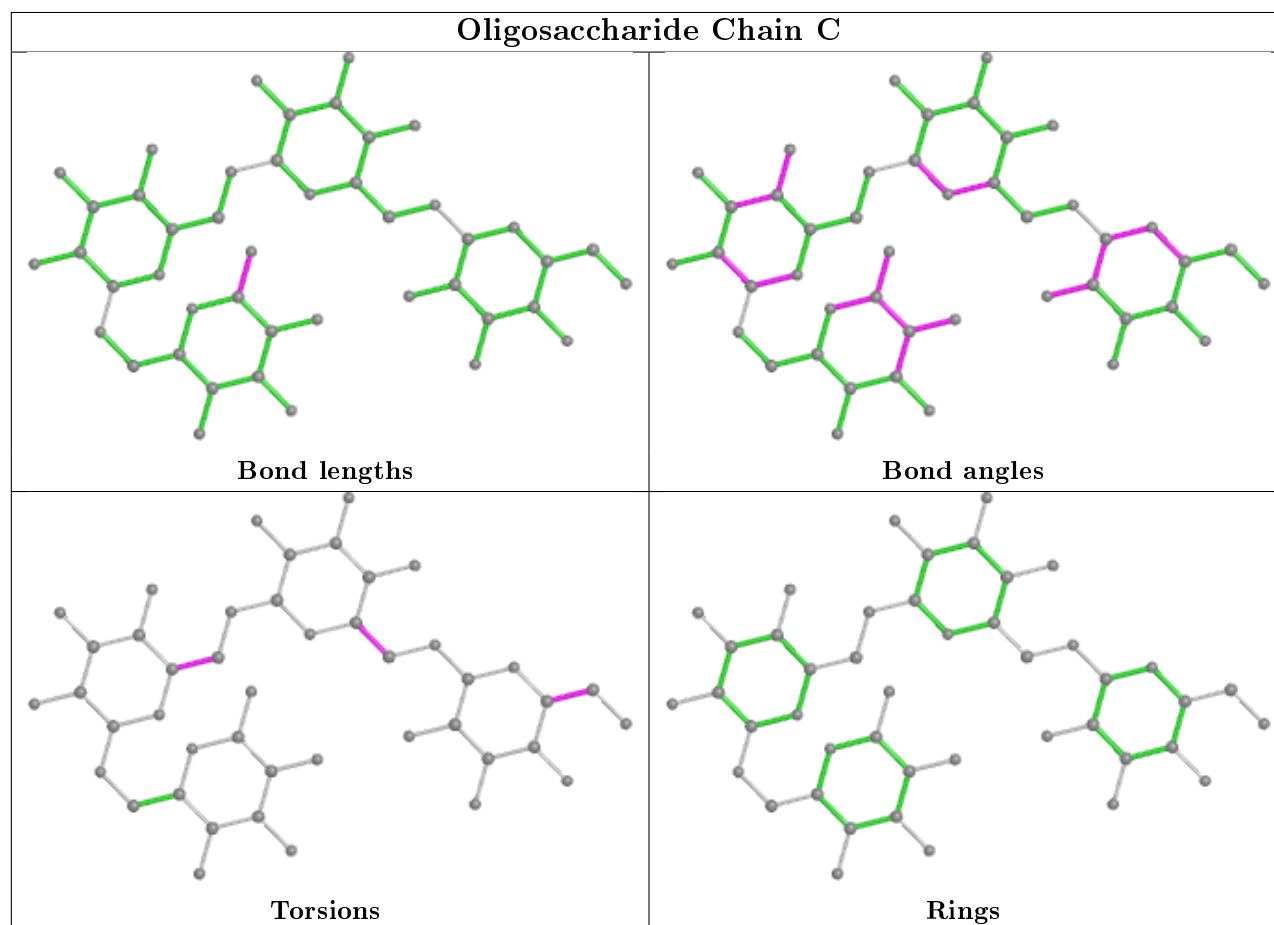
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	1	0
2	C	3	GLC	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4	GLC	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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1265/1290 (98%)	-0.33	0 100 100	41, 57, 87, 106	0
1	B	1144/1290 (88%)	-0.31	2 (0%) 95 93	46, 65, 88, 110	0
All	All	2409/2580 (93%)	-0.32	2 (0%) 95 94	41, 61, 88, 110	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	313	ASN	2.6
1	B	302	VAL	2.5

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

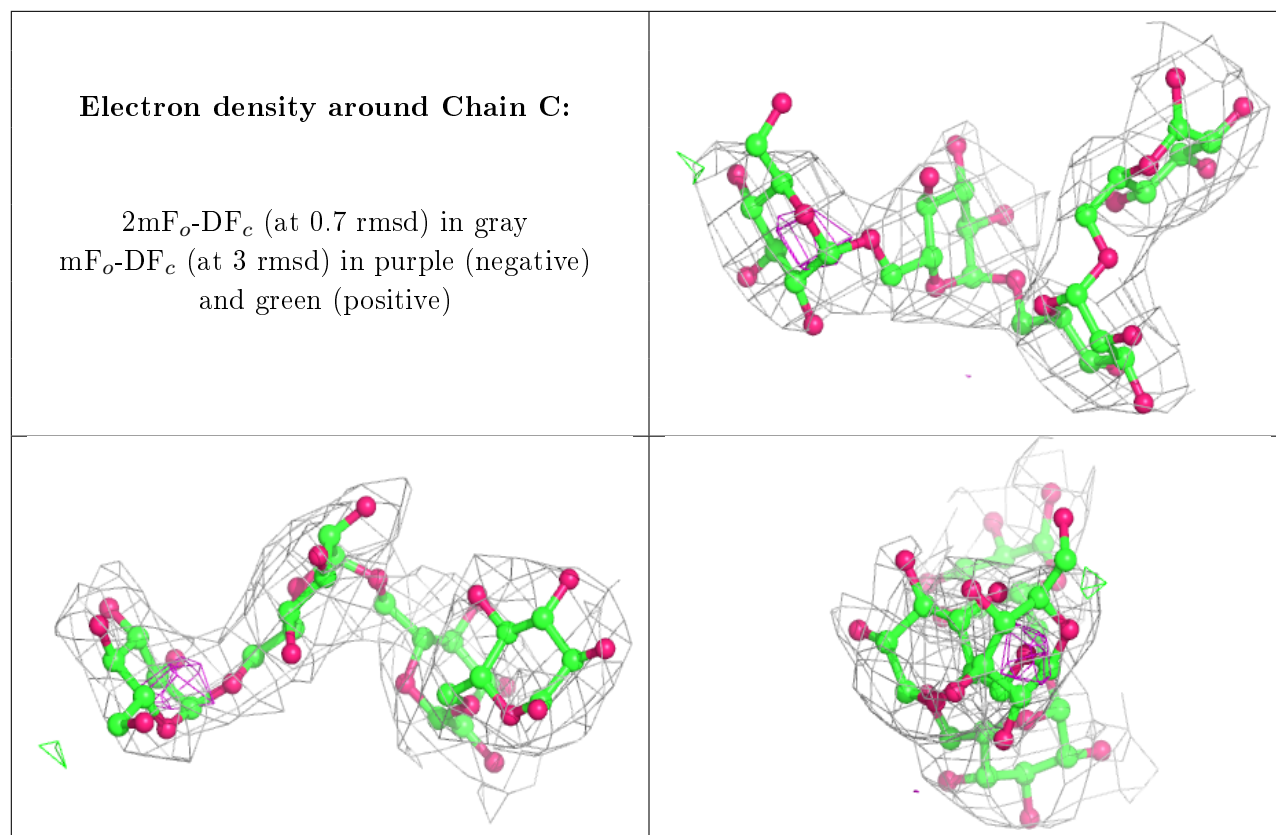
There are no non-standard protein/DNA/RNA residues in this entry.

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	GLC	C	4	11/12	0.77	0.47	76,81,85,85	0
2	GLC	C	3	11/12	0.84	0.30	69,74,76,81	0
2	GLC	C	1	12/12	0.92	0.18	59,61,63,64	0
2	GLC	C	2	11/12	0.95	0.24	60,64,69,72	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](#)

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	PR	A	1504	1/1	0.89	0.14	82,82,82,82	1
4	PR	A	1503	1/1	0.94	0.27	69,69,69,69	1
4	PR	B	1504	1/1	0.96	0.13	84,84,84,84	1
4	PR	B	1503	1/1	0.96	0.26	59,59,59,59	1
4	PR	B	1502	1/1	0.96	0.26	54,54,54,54	1
3	CA	A	1501	1/1	0.98	0.11	41,41,41,41	0
3	CA	B	1501	1/1	0.99	0.06	44,44,44,44	0
4	PR	A	1502	1/1	0.99	0.19	79,79,79,79	1

6.5 Other polymers [i](#)

There are no such residues in this entry.