



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:03 PM UTC

PDB ID : 2FHF / pdb_00002fhf
Title : Crystal Structure Analysis of Klebsiella pneumoniae pullulanase complexed with maltotetraose
Authors : Mikami, B.; Iwamoto, H.; Katsuya, Y.; Yoon, H.-J.; Demirkan-Sarikaya, E.; Malle, D.
Deposited on : 2005-12-23
Resolution : 1.65 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

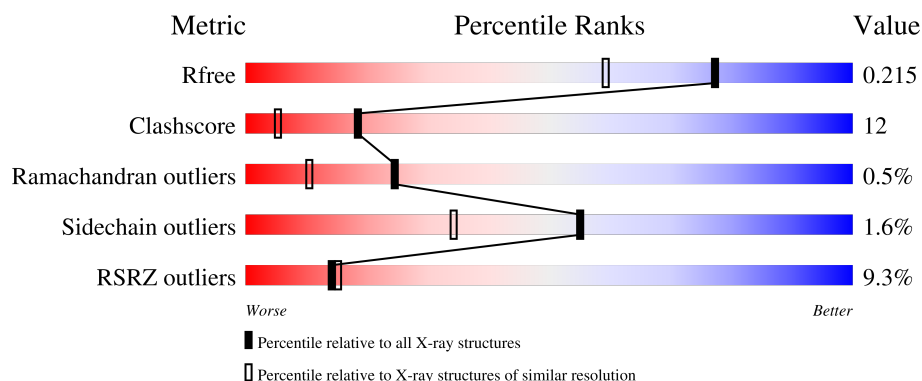
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	1083	9% 76% 19% • •
2	B	4	75% 25%
2	C	4	50% 50%
3	D	2	50% 50%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9441 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 Alpha-dextrin endo-1,6-alpha-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1052	Total	C	N	O	S	0	18	0
			8126	5081	1382	1636	27			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	680	LEU	GLY	conflict	UNP W9BQ28

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



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

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total 5	Ca 5	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1197	Total 1197	O 1197	0	0

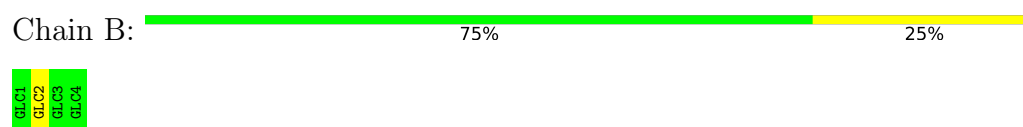
3 Residue-property plots

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: Alpha-dextrin endo-1,6-alpha-glucosidase



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



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

Chain C:  50% 50%

GLC1
GLC2
GLC3
GLC4

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:  50% 50%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	150.16Å 60.51Å 134.66Å 90.00° 111.87° 90.00°	Depositor
Resolution (Å)	14.98 – 1.65 14.98 – 1.65	Depositor EDS
% Data completeness (in resolution range)	85.7 (14.98-1.65) 85.6 (14.98-1.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.66 (at 1.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.178 , 0.206 0.188 , 0.215	Depositor DCC
R_{free} test set	11574 reflections (9.25%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.5	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	9441	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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: 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	0.35	0/8363	0.90	35/11388 (0.3%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	644	CYS	N-CA-CB	-9.45	102.80	111.59
1	A	555	TYR	N-CA-C	7.72	121.50	109.07
1	A	484	THR	N-CA-C	7.47	121.81	113.21
1	A	652	ARG	N-CA-C	7.19	119.74	111.11
1	A	289	ALA	N-CA-C	6.99	119.50	111.11
1	A	426	ILE	N-CA-C	6.95	119.43	108.87
1	A	480	PHE	N-CA-C	-6.78	99.86	110.36
1	A	845	LYS	N-CA-C	6.73	121.78	113.50
1	A	931	LYS	N-CA-C	6.63	119.06	111.11
1	A	644	CYS	N-CA-C	6.59	118.02	108.34
1	A	562	PHE	N-CA-C	-6.33	104.51	112.93
1	A	114	ILE	N-CA-C	6.25	114.08	107.76
1	A	280	THR	N-CA-C	5.79	120.04	112.92
1	A	820	ALA	N-CA-C	5.75	119.03	110.52
1	A	904	SER	N-CA-C	-5.75	106.04	113.16
1	A	718	ILE	N-CA-C	5.70	116.27	110.05
1	A	297	TYR	N-CA-C	5.57	118.32	109.81
1	A	532	LYS	N-CA-C	5.48	116.93	111.07
1	A	312	ALA	CA-C-N	5.40	125.07	119.56
1	A	312	ALA	C-N-CA	5.40	125.07	119.56
1	A	1019	SER	N-CA-C	-5.40	101.67	110.20
1	A	891	SER	N-CA-C	5.39	118.97	112.93
1	A	578	THR	N-CA-C	5.29	119.36	113.01
1	A	471	VAL	N-CA-C	-5.26	102.73	109.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	883	ARG	N-CA-C	5.26	117.97	110.24
1	A	435	SER	N-CA-C	5.23	121.46	113.61
1	A	764	VAL	N-CA-C	5.20	118.41	111.44
1	A	520	SER	N-CA-C	-5.20	103.17	110.35
1	A	278	VAL	N-CA-C	5.20	115.44	108.17
1	A	182	VAL	N-CA-C	5.16	118.35	111.44
1	A	722	ILE	N-CA-C	5.07	115.22	110.30
1	A	429	SER	N-CA-C	5.06	117.55	109.81
1	A	564	TYR	N-CA-C	5.04	118.69	112.54
1	A	905	LEU	N-CA-C	5.01	118.66	112.24
1	A	897	TRP	N-CA-C	5.01	116.43	110.97

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	8126	0	7830	193	0
2	B	45	0	39	2	0
2	C	45	0	39	4	0
3	D	23	0	21	2	0
4	A	5	0	0	0	0
5	A	1197	0	0	29	0
All	All	9441	0	7929	195	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 12.

All (195) 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:740:VAL:HA	1:A:758:VAL:HG12	1.44	0.97
1:A:81:ASN:HB2	1:A:86:ASP:HA	1.54	0.89
1:A:606:ASN:HD21	1:A:607:HIS:HD2	1.20	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:978:ASN:HD21	1:A:984:GLN:H	1.20	0.88
2:B:2:GLC:H62	2:C:1:GLC:O1	1.77	0.85
1:A:722:ILE:HD11	5:A:1973:HOH:O	1.78	0.83
1:A:677:ASP:OD1	2:C:1:GLC:H1	1.80	0.82
1:A:81:ASN:HB3	1:A:88:LEU:HD13	1.62	0.80
1:A:680:LEU:HD11	5:A:1539:HOH:O	1.82	0.80
1:A:195:LYS:HG2	1:A:265:ALA:HB1	1.63	0.79
1:A:977:ARG:NH1	1:A:1024:ASP:HB3	1.98	0.79
1:A:230:GLN:HG2	1:A:234:MET:HE2	1.64	0.78
1:A:722:ILE:HG12	5:A:1660:HOH:O	1.84	0.78
1:A:79:LEU:HD11	1:A:114:ILE:HD12	1.66	0.77
1:A:606:ASN:ND2	1:A:607:HIS:HD2	1.82	0.77
1:A:516:GLY:HA3	5:A:1195:HOH:O	1.85	0.77
1:A:324:SER:HB3	1:A:330[B]:ILE:HD11	1.68	0.76
1:A:157:ASN:HD21	1:A:159:ALA:HB3	1.50	0.75
1:A:229:ASN:ND2	1:A:232:VAL:H	1.88	0.72
1:A:501:ARG:O	1:A:504:GLU:HG2	1.89	0.72
1:A:560:ASP:HB3	1:A:609:ASN:ND2	2.03	0.72
1:A:680:LEU:HG	1:A:710:SER:HB3	1.73	0.71
1:A:560:ASP:HB3	1:A:609:ASN:HD22	1.57	0.70
1:A:376:VAL:HB	1:A:623[B]:ILE:HD11	1.74	0.69
1:A:88:LEU:HD12	1:A:88:LEU:H	1.58	0.68
1:A:76:ASN:CG	1:A:98:VAL:HG23	2.19	0.68
1:A:229:ASN:HD21	1:A:232:VAL:HG23	1.57	0.68
1:A:740:VAL:HA	1:A:758:VAL:CG1	2.20	0.67
1:A:1030:LEU:HG	1:A:1066:LEU:HB3	1.76	0.67
1:A:505[A]:VAL:HG22	5:A:1366:HOH:O	1.95	0.66
1:A:107:LYS:HE3	1:A:107:LYS:H	1.61	0.66
1:A:627:TYR:O	1:A:651:HIS:HD2	1.79	0.66
1:A:157:ASN:ND2	1:A:159:ALA:HB3	2.11	0.65
1:A:84:THR:HG21	1:A:140:ARG:HH11	1.61	0.65
1:A:977:ARG:HH11	1:A:1024:ASP:HB3	1.60	0.65
1:A:680:LEU:HD13	1:A:708:TRP:HB2	1.78	0.64
1:A:59:VAL:HB	1:A:154:ILE:HG12	1.81	0.62
1:A:1005:ARG:HB3	1:A:1005:ARG:NH1	2.14	0.62
1:A:677:ASP:OD1	2:C:1:GLC:C1	2.47	0.62
1:A:127:VAL:HB	1:A:135:ILE:HD13	1.81	0.62
1:A:127:VAL:HG12	1:A:134:LEU:HD12	1.81	0.61
1:A:83:GLU:HG3	1:A:84:THR:N	2.16	0.60
1:A:229:ASN:HD21	1:A:232:VAL:CG2	2.15	0.60
1:A:246:LYS:HD3	1:A:247:LEU:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:LYS:C	1:A:505[A]:VAL:HG21	2.28	0.59
1:A:682:HIS:HD2	1:A:686:GLN:NE2	2.01	0.58
1:A:134:LEU:C	1:A:135:ILE:HD12	2.28	0.58
1:A:495:ILE:HA	1:A:524:VAL:HB	1.86	0.58
1:A:1028:THR:HG22	1:A:1068:LEU:HD21	1.86	0.58
1:A:129:ASP:HB2	1:A:134:LEU:HD21	1.86	0.57
1:A:84:THR:HG21	1:A:140:ARG:NH1	2.19	0.57
1:A:816:PRO:HG2	5:A:1973:HOH:O	2.03	0.57
1:A:680:LEU:HD23	1:A:680:LEU:N	2.19	0.57
1:A:606:ASN:ND2	1:A:607:HIS:CD2	2.68	0.57
1:A:98:VAL:HG22	1:A:98:VAL:O	2.05	0.57
1:A:157:ASN:HB3	5:A:2296:HOH:O	2.05	0.57
1:A:722:ILE:HD12	5:A:1974:HOH:O	2.05	0.57
1:A:56:ILE:CG2	1:A:112:TRP:HB2	2.35	0.56
1:A:83:GLU:HG3	1:A:84:THR:H	1.69	0.56
1:A:680:LEU:HD23	1:A:680:LEU:H	1.69	0.56
1:A:682:HIS:HD2	1:A:686:GLN:HE22	1.53	0.56
1:A:1049:GLN:HG3	1:A:1057:THR:HB	1.87	0.56
1:A:126:ILE:HD11	1:A:138:ASP:OD1	2.05	0.56
1:A:229:ASN:HD21	1:A:232:VAL:H	1.51	0.56
1:A:642:THR:O	1:A:643:CYS:HB3	2.06	0.56
1:A:466:LEU:HD23	1:A:950:GLN:HE21	1.71	0.55
1:A:108:TYR:HB3	5:A:2279:HOH:O	2.05	0.55
1:A:56:ILE:HG22	1:A:112:TRP:HB2	1.87	0.55
1:A:198:VAL:HG12	1:A:223:LEU:HD12	1.89	0.55
1:A:774:ASP:HB3	5:A:2292:HOH:O	2.06	0.54
1:A:87:ALA:O	1:A:117:THR:HG22	2.07	0.54
1:A:85:CYS:HB3	1:A:122:CYS:C	2.32	0.54
1:A:254:ASP:HB2	5:A:2122:HOH:O	2.07	0.54
1:A:387:ASN:ND2	1:A:487:GLU:H	2.05	0.54
1:A:106:ASP:HB2	1:A:107:LYS:NZ	2.23	0.54
1:A:1005:ARG:HB3	1:A:1005:ARG:HH11	1.73	0.54
1:A:496:GLN:NE2	1:A:496:GLN:H	2.05	0.53
1:A:1039:ALA:HB3	1:A:1043:SER:HB2	1.90	0.53
1:A:430:HIS:CD2	1:A:432:ARG:H	2.27	0.53
1:A:680:LEU:HD12	1:A:709:ASP:C	2.35	0.53
1:A:946:THR:O	1:A:950:GLN:HG3	2.09	0.53
1:A:465:GLN:HG3	1:A:950:GLN:HE22	1.75	0.52
1:A:229:ASN:C	1:A:229:ASN:HD22	2.17	0.52
1:A:678:LEU:C	1:A:680:LEU:HD23	2.35	0.52
1:A:680:LEU:HD12	1:A:710:SER:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ALA:HB3	5:A:2256:HOH:O	2.10	0.51
1:A:79:LEU:HB3	1:A:88:LEU:HD23	1.91	0.51
1:A:43:PRO:HB3	1:A:108:TYR:CG	2.46	0.51
1:A:682:HIS:CD2	1:A:686:GLN:HE22	2.29	0.51
1:A:40:VAL:HG13	1:A:41:ALA:N	2.26	0.51
1:A:229:ASN:HD21	1:A:232:VAL:CB	2.24	0.50
1:A:680:LEU:H	1:A:680:LEU:CD2	2.25	0.50
1:A:107:LYS:H	1:A:107:LYS:CE	2.24	0.50
1:A:272:LEU:C	1:A:272:LEU:HD23	2.36	0.50
1:A:642:THR:O	1:A:643:CYS:CB	2.60	0.50
1:A:54:ALA:HB2	1:A:116:LEU:HD11	1.94	0.49
1:A:38:PRO:O	1:A:169:ARG:HD3	2.12	0.49
1:A:85:CYS:HB3	1:A:122:CYS:O	2.12	0.49
1:A:61:ILE:O	1:A:61:ILE:HG13	2.12	0.49
1:A:117:THR:HG23	1:A:118:LYS:HG2	1.94	0.49
1:A:80:TRP:O	1:A:124:ASN:HB2	2.12	0.49
1:A:523:THR:OG1	1:A:526:GLU:HG3	2.13	0.48
1:A:76:ASN:ND2	1:A:128:ARG:HH21	2.11	0.48
1:A:204:HIS:HE1	5:A:1154:HOH:O	1.97	0.48
1:A:81:ASN:ND2	1:A:91:PRO:HG3	2.29	0.48
1:A:182:VAL:HG21	1:A:232:VAL:HG11	1.96	0.48
1:A:978:ASN:ND2	1:A:984:GLN:H	2.01	0.48
1:A:88:LEU:HD12	1:A:88:LEU:N	2.26	0.48
1:A:411[A]:MET:HE2	1:A:672:ASP:OD1	2.14	0.48
1:A:107:LYS:HD3	1:A:108:TYR:CE2	2.48	0.47
1:A:1028:THR:HG22	1:A:1068:LEU:CD2	2.43	0.47
1:A:88:LEU:HD11	1:A:123:ILE:HA	1.95	0.47
1:A:1066:LEU:HD22	1:A:1066:LEU:N	2.28	0.47
1:A:740:VAL:HG22	1:A:758:VAL:CG1	2.45	0.47
1:A:657:LEU:C	1:A:657:LEU:HD23	2.40	0.47
1:A:176:LEU:HA	5:A:1715:HOH:O	2.14	0.47
1:A:72:TYR:CZ	1:A:110:PRO:HD3	2.50	0.47
1:A:157:ASN:O	1:A:158:SER:HB3	2.15	0.47
1:A:430:HIS:HD2	1:A:433:ASP:H	1.62	0.47
1:A:461[A]:GLN:NE2	5:A:1931:HOH:O	2.48	0.47
1:A:1014:ASN:HD21	1:A:1020:ARG:HH11	1.61	0.46
1:A:778:ARG:HD2	5:A:1519:HOH:O	2.14	0.46
1:A:253:VAL:HG23	1:A:254:ASP:N	2.29	0.46
1:A:105:SER:HB3	5:A:2282:HOH:O	2.14	0.46
1:A:260:GLU:HB2	1:A:364:TYR:CE1	2.51	0.46
1:A:107:LYS:HD3	1:A:108:TYR:CD2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:HIS:H	1:A:238:HIS:CD2	2.33	0.46
1:A:92:VAL:HG21	1:A:97:ASP:OD2	2.16	0.46
1:A:797:LEU:HD12	1:A:797:LEU:C	2.40	0.46
1:A:147:THR:HG23	1:A:148:ASP:N	2.31	0.46
1:A:95:TRP:CZ3	3:D:1:GLC:H2	2.51	0.45
1:A:465:GLN:HG3	1:A:950:GLN:NE2	2.30	0.45
1:A:59:VAL:HG13	5:A:2279:HOH:O	2.16	0.45
1:A:1049:GLN:CG	1:A:1057:THR:HB	2.47	0.45
1:A:199:ARG:HD3	1:A:220:TYR:CE2	2.51	0.45
1:A:377:THR:OG1	1:A:563:HIS:HE1	1.99	0.45
1:A:252:ASN:C	1:A:252:ASN:HD22	2.23	0.45
1:A:525:GLU:HG2	5:A:2178:HOH:O	2.15	0.45
1:A:238:HIS:HE1	5:A:1433:HOH:O	1.99	0.45
1:A:383:SER:HB3	1:A:392:GLN:HB3	1.98	0.45
1:A:80:TRP:CZ2	1:A:124:ASN:HB3	2.52	0.44
1:A:366:PRO:HB2	1:A:626:TRP:CE2	2.53	0.44
1:A:491:LYS:O	1:A:505[A]:VAL:HG21	2.17	0.44
1:A:199:ARG:HD3	1:A:220:TYR:CD2	2.52	0.44
1:A:643:CYS:SG	1:A:644:CYS:N	2.91	0.44
1:A:798[B]:ILE:HD13	5:A:1401:HOH:O	2.18	0.44
1:A:449:TYR:CZ	1:A:571:TYR:HB2	2.53	0.43
1:A:785:ARG:NH2	5:A:1359:HOH:O	2.50	0.43
1:A:886:SER:O	1:A:887:PHE:HB2	2.19	0.43
1:A:40:VAL:HG13	1:A:41:ALA:H	1.83	0.43
1:A:740:VAL:HG22	1:A:758:VAL:HG11	2.00	0.43
1:A:675:ARG:HH21	1:A:832:LYS:HZ1	1.67	0.43
1:A:229:ASN:ND2	1:A:229:ASN:C	2.77	0.43
1:A:229:ASN:ND2	1:A:232:VAL:HB	2.33	0.43
1:A:269:ASP:OD1	1:A:271:ILE:HG13	2.18	0.43
1:A:36:ARG:HG2	1:A:36:ARG:HH11	1.84	0.43
1:A:56:ILE:HG21	1:A:77:LEU:HD11	2.00	0.43
1:A:135:ILE:HD12	1:A:135:ILE:N	2.34	0.43
1:A:272:LEU:HD23	1:A:273:SER:N	2.33	0.43
2:B:2:GLC:H62	2:C:1:GLC:HO1	1.81	0.43
1:A:106:ASP:HB2	1:A:107:LYS:HZ1	1.84	0.42
1:A:413:HIS:HD2	5:A:1172:HOH:O	2.02	0.42
1:A:418:LYS:HD2	1:A:966[B]:ASP:OD1	2.19	0.42
1:A:52:ARG:HH11	1:A:52:ARG:HG3	1.84	0.42
1:A:187:LEU:C	1:A:187:LEU:HD13	2.44	0.42
1:A:56:ILE:HG23	1:A:112:TRP:HE3	1.84	0.42
1:A:75:LYS:O	1:A:102:PRO:HD3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:SER:H	1:A:53:GLN:NE2	2.18	0.42
1:A:231:GLN:OE1	1:A:231:GLN:N	2.51	0.42
1:A:465:GLN:CG	1:A:950:GLN:HE22	2.33	0.42
1:A:606:ASN:HD21	1:A:607:HIS:CD2	2.13	0.41
1:A:682:HIS:CD2	1:A:686:GLN:NE2	2.85	0.41
1:A:108:TYR:HB2	1:A:164:ARG:NH1	2.34	0.41
1:A:95:TRP:CH2	3:D:1:GLC:H2	2.56	0.41
1:A:1005:ARG:HH11	1:A:1005:ARG:CB	2.34	0.41
1:A:481:ASP:OD2	1:A:563:HIS:HD2	2.03	0.41
1:A:543:GLN:NE2	1:A:917:SER:H	2.19	0.41
1:A:939:GLU:OE1	1:A:943:LYS:HE3	2.21	0.41
1:A:110:PRO:HB3	5:A:2282:HOH:O	2.20	0.41
1:A:58:LEU:HB2	5:A:2283:HOH:O	2.21	0.41
1:A:494:ASP:HB3	1:A:496:GLN:HE22	1.85	0.41
1:A:875:ASP:OD1	1:A:879:SER:HB2	2.21	0.41
1:A:387:ASN:HD22	1:A:487:GLU:HG3	1.86	0.41
1:A:156:GLY:HA2	5:A:2222:HOH:O	2.20	0.41
1:A:226:THR:HB	1:A:246:LYS:HB2	2.03	0.41
1:A:929:ARG:HD2	5:A:1584:HOH:O	2.20	0.41
1:A:173:GLY:HA2	5:A:2189:HOH:O	2.21	0.41
1:A:848:GLN:HG3	5:A:1506:HOH:O	2.21	0.41
1:A:254:ASP:O	1:A:258:GLN:HG2	2.22	0.40
1:A:388:SER:HB2	1:A:562:PHE:CE1	2.55	0.40
1:A:555:TYR:CD1	1:A:555:TYR:C	2.99	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	1068/1083 (99%)	1036 (97%)	27 (2%)	5 (0%)	24 10

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	478	PRO
1	A	643	CYS
1	A	136	ASP
1	A	94	ASP
1	A	137	SER

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	884/891 (99%)	870 (98%)	14 (2%)	55 34

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	ILE
1	A	75	LYS
1	A	107	LYS
1	A	195	LYS
1	A	229	ASN
1	A	374	TYR
1	A	387	ASN
1	A	478	PRO
1	A	482	LEU
1	A	496	GLN
1	A	543	GLN
1	A	859	GLN
1	A	886	SER
1	A	1066	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN

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Mol	Chain	Res	Type
1	A	57	HIS
1	A	81	ASN
1	A	157	ASN
1	A	194	ASN
1	A	204	HIS
1	A	213	ASN
1	A	229	ASN
1	A	230	GLN
1	A	238	HIS
1	A	252	ASN
1	A	279	GLN
1	A	316	GLN
1	A	317	GLN
1	A	387	ASN
1	A	392	GLN
1	A	413	HIS
1	A	430	HIS
1	A	439	GLN
1	A	455	GLN
1	A	458	ASN
1	A	465	GLN
1	A	496	GLN
1	A	533	GLN
1	A	534	ASN
1	A	541	GLN
1	A	543	GLN
1	A	551	GLN
1	A	563	HIS
1	A	606	ASN
1	A	607	HIS
1	A	609	ASN
1	A	629	GLN
1	A	651	HIS
1	A	682	HIS
1	A	686	GLN
1	A	711	ASN
1	A	712	GLN
1	A	835	ASN
1	A	848	GLN
1	A	859	GLN
1	A	906	GLN
1	A	911	ASN

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Mol	Chain	Res	Type
1	A	950	GLN
1	A	978	ASN
1	A	984	GLN
1	A	997	GLN
1	A	1014	ASN
1	A	1023	GLN
1	A	1031	GLN
1	A	1037	GLN
1	A	1049	GLN
1	A	1074	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 ⓘ

10 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	B	1	2	12,12,12	0.48	0	17,17,17	0.51	0
2	GLC	B	2	2	11,11,12	0.50	0	15,15,17	0.65	0
2	GLC	B	3	2	11,11,12	0.52	0	15,15,17	0.55	0
2	GLC	B	4	2	11,11,12	0.47	0	15,15,17	0.59	0
2	GLC	C	1	2	12,12,12	0.42	0	17,17,17	0.58	0
2	GLC	C	2	2	11,11,12	0.52	0	15,15,17	0.69	1 (6%)
2	GLC	C	3	2	11,11,12	0.49	0	15,15,17	0.61	0
2	GLC	C	4	2	11,11,12	0.44	0	15,15,17	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	D	1	3	12,12,12	0.42	0	17,17,17	0.32	0
3	GLC	D	2	3	11,11,12	0.47	0	15,15,17	0.57	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
2	GLC	B	1	2	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	2/2/19/22	0/1/1/1
2	GLC	B	3	2	-	2/2/19/22	0/1/1/1
2	GLC	B	4	2	-	0/2/19/22	0/1/1/1
2	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	C	3	2	-	0/2/19/22	0/1/1/1
2	GLC	C	4	2	-	0/2/19/22	0/1/1/1
3	GLC	D	1	3	-	0/2/22/22	0/1/1/1
3	GLC	D	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	GLC	C1-O5-C5	2.42	115.42	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

There are no ring outliers.

3 monomers are involved in 6 short contacts:

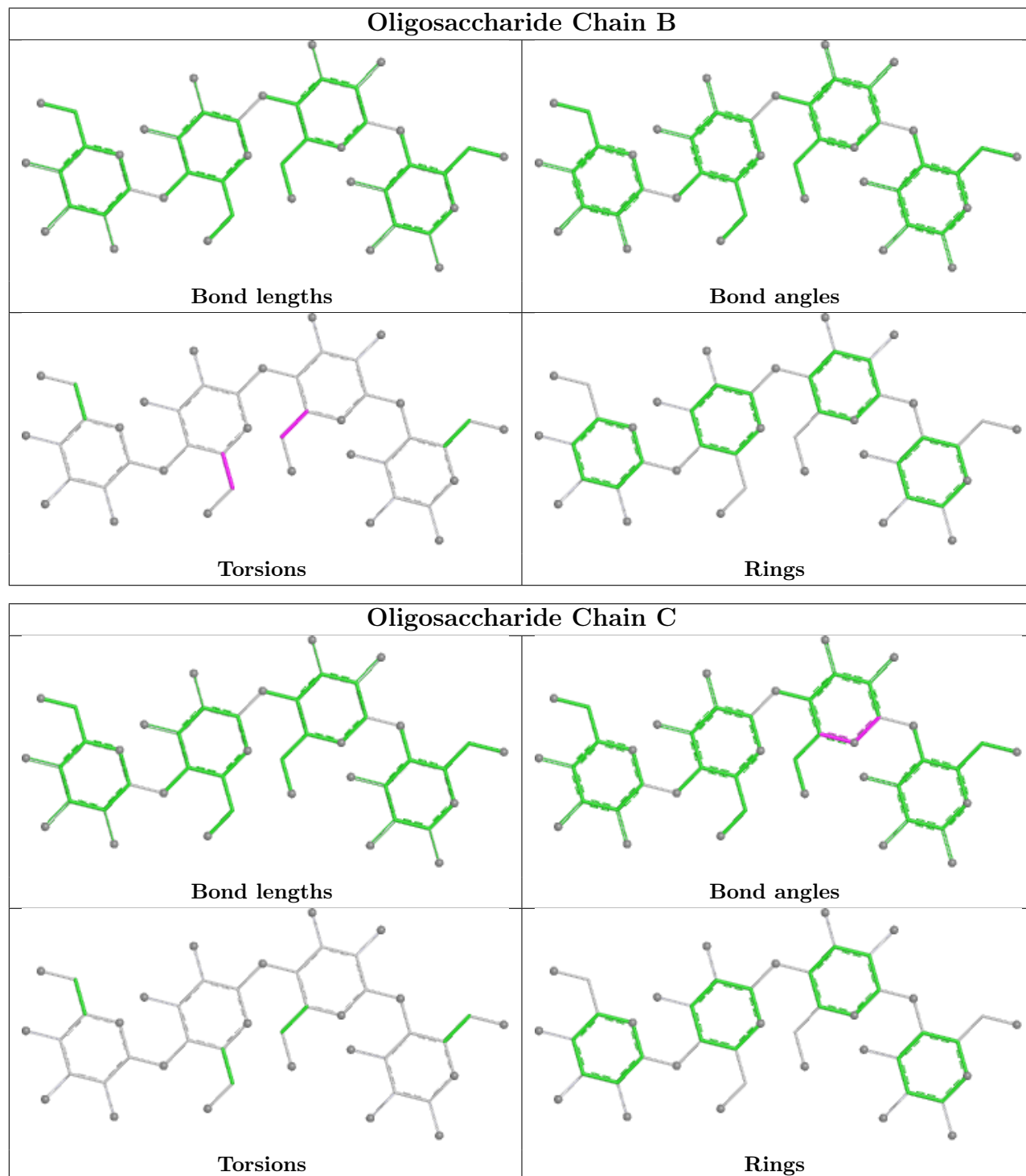
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	4	0

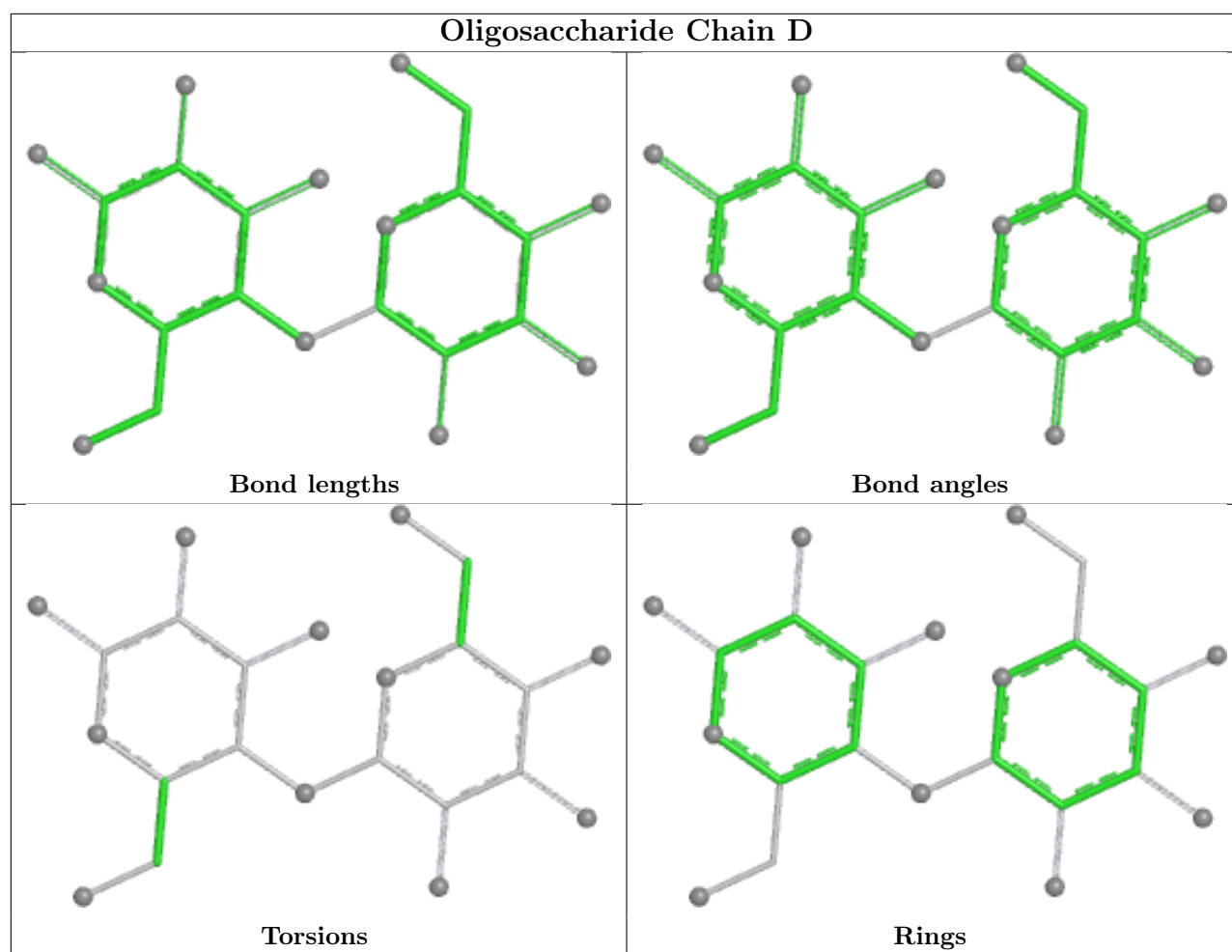
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	GLC	2	0
2	B	2	GLC	2	0

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





5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 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 [i](#)

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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1052/1083 (97%)	0.11	98 (9%) 14 15	12, 20, 50, 63	142 (13%)

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	VAL	5.8
1	A	303	ASP	5.3
1	A	93	ALA	4.7
1	A	81	ASN	4.7
1	A	72	TYR	4.7
1	A	135	ILE	4.5
1	A	56	ILE	4.4
1	A	62	ALA	4.0
1	A	158	SER	3.9
1	A	95	TRP	3.8
1	A	78	TYR	3.8
1	A	126	ILE	3.7
1	A	117	THR	3.6
1	A	129	ASP	3.6
1	A	127	VAL	3.6
1	A	134	LEU	3.6
1	A	79	LEU	3.5
1	A	154	ILE	3.5
1	A	42	VAL	3.5
1	A	92	VAL	3.4
1	A	131	THR	3.4
1	A	168	PHE	3.3
1	A	172	PHE	3.3
1	A	61	ILE	3.3
1	A	123	ILE	3.3
1	A	88	LEU	3.3
1	A	77	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	161	TYR	3.2
1	A	35	VAL	3.2
1	A	136	ASP	3.1
1	A	157	ASN	3.1
1	A	84	THR	3.1
1	A	147	THR	3.0
1	A	59	VAL	3.0
1	A	87	ALA	3.0
1	A	159	ALA	3.0
1	A	112	TRP	2.9
1	A	34	VAL	2.9
1	A	98	VAL	2.9
1	A	90	ALA	2.9
1	A	130	GLY	2.9
1	A	99	SER	2.9
1	A	120	SER	2.9
1	A	74	THR	2.8
1	A	73	ALA	2.8
1	A	165	ALA	2.8
1	A	108	TYR	2.8
1	A	33	VAL	2.8
1	A	47	VAL	2.7
1	A	247	LEU	2.7
1	A	58	LEU	2.7
1	A	69	PRO	2.7
1	A	80	TRP	2.7
1	A	83	GLU	2.6
1	A	68	THR	2.6
1	A	109	GLY	2.6
1	A	113	VAL	2.6
1	A	208	VAL	2.6
1	A	138	ASP	2.6
1	A	302	THR	2.6
1	A	156	GLY	2.5
1	A	146	PHE	2.5
1	A	212	SER	2.5
1	A	153	VAL	2.5
1	A	64	ILE	2.5
1	A	746	PHE	2.5
1	A	171	ALA	2.5
1	A	144	SER	2.5
1	A	55	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	268	SER	2.4
1	A	85	CYS	2.4
1	A	70	ALA	2.3
1	A	76	ASN	2.3
1	A	116	LEU	2.3
1	A	142	SER	2.3
1	A	125	VAL	2.3
1	A	167	ALA	2.2
1	A	82	ASN	2.2
1	A	67	SER	2.2
1	A	163	SER	2.2
1	A	32	ASP	2.2
1	A	1030	LEU	2.2
1	A	151	VAL	2.1
1	A	41	ALA	2.1
1	A	155	ALA	2.1
1	A	104	GLY	2.1
1	A	139	LEU	2.1
1	A	94	ASP	2.1
1	A	65	THR	2.1
1	A	102	PRO	2.1
1	A	248	PRO	2.1
1	A	143	PHE	2.1
1	A	271	ILE	2.1
1	A	128	ARG	2.1
1	A	114	ILE	2.0
1	A	197	ILE	2.0
1	A	71	ASP	2.0
1	A	218	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

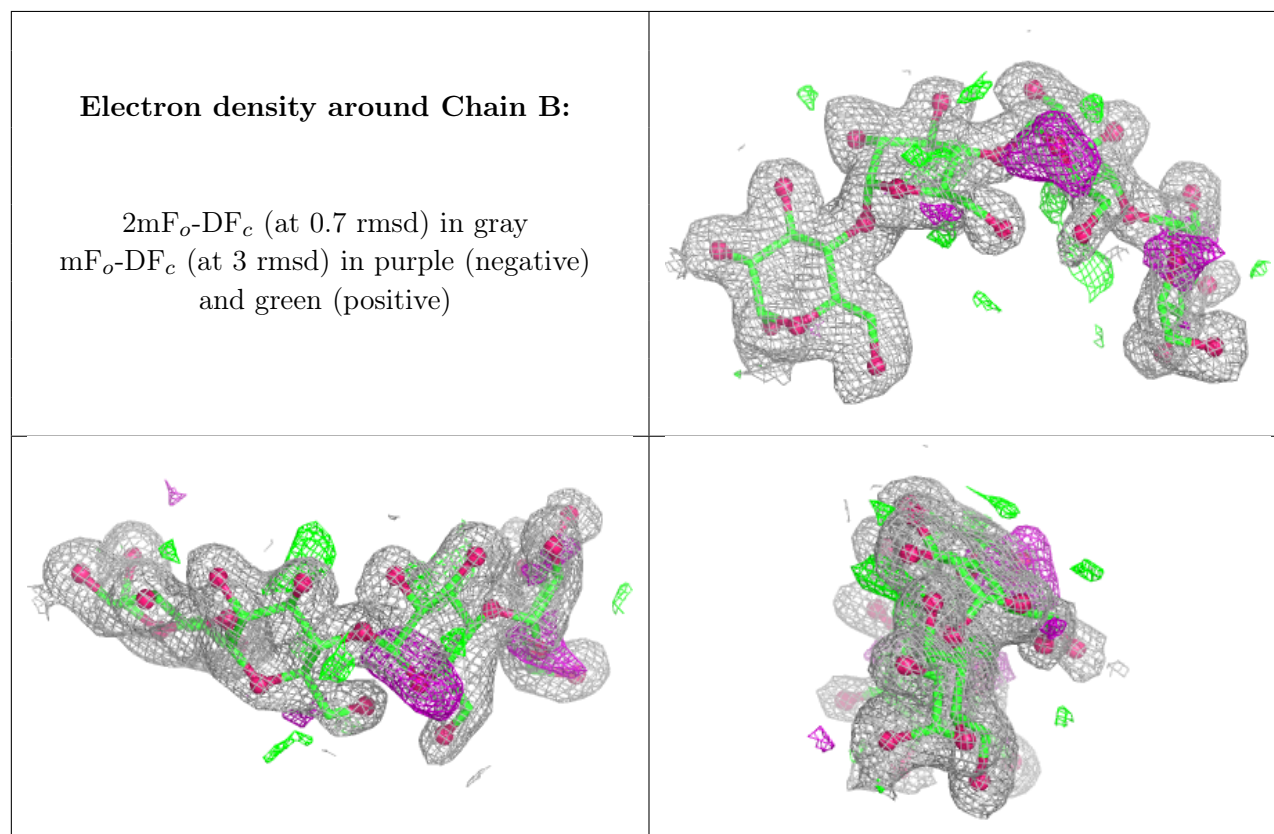
There are no non-standard protein/DNA/RNA residues in this entry.

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.

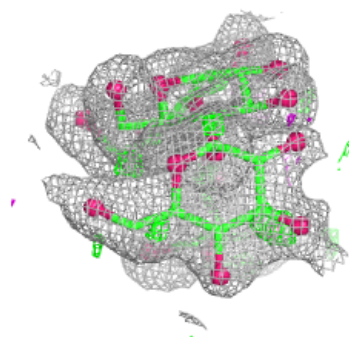
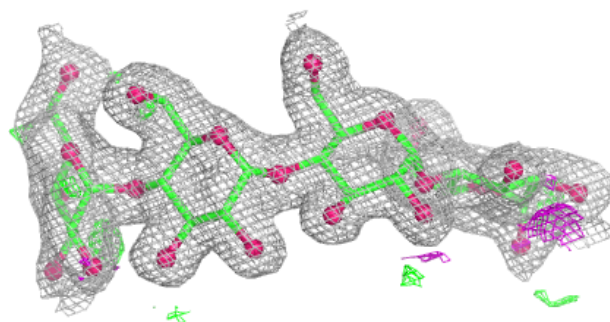
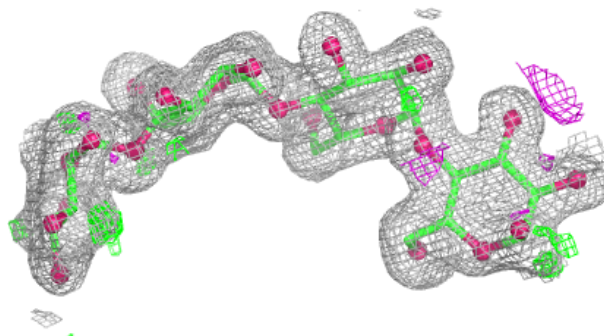
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GLC	D	1	12/12	0.72	0.12	60,61,62,62	0
2	GLC	C	4	11/12	0.74	0.12	32,35,38,40	0
3	GLC	D	2	11/12	0.76	0.10	62,63,63,63	0
2	GLC	B	4	11/12	0.83	0.10	33,35,36,36	0
2	GLC	B	3	11/12	0.83	0.10	27,31,33,34	0
2	GLC	B	2	11/12	0.86	0.09	26,30,33,35	0
2	GLC	C	1	12/12	0.92	0.08	18,23,26,31	0
2	GLC	C	3	11/12	0.94	0.06	20,24,28,29	0
2	GLC	B	1	12/12	0.95	0.07	19,22,23,23	0
2	GLC	C	2	11/12	0.97	0.04	14,15,17,17	0

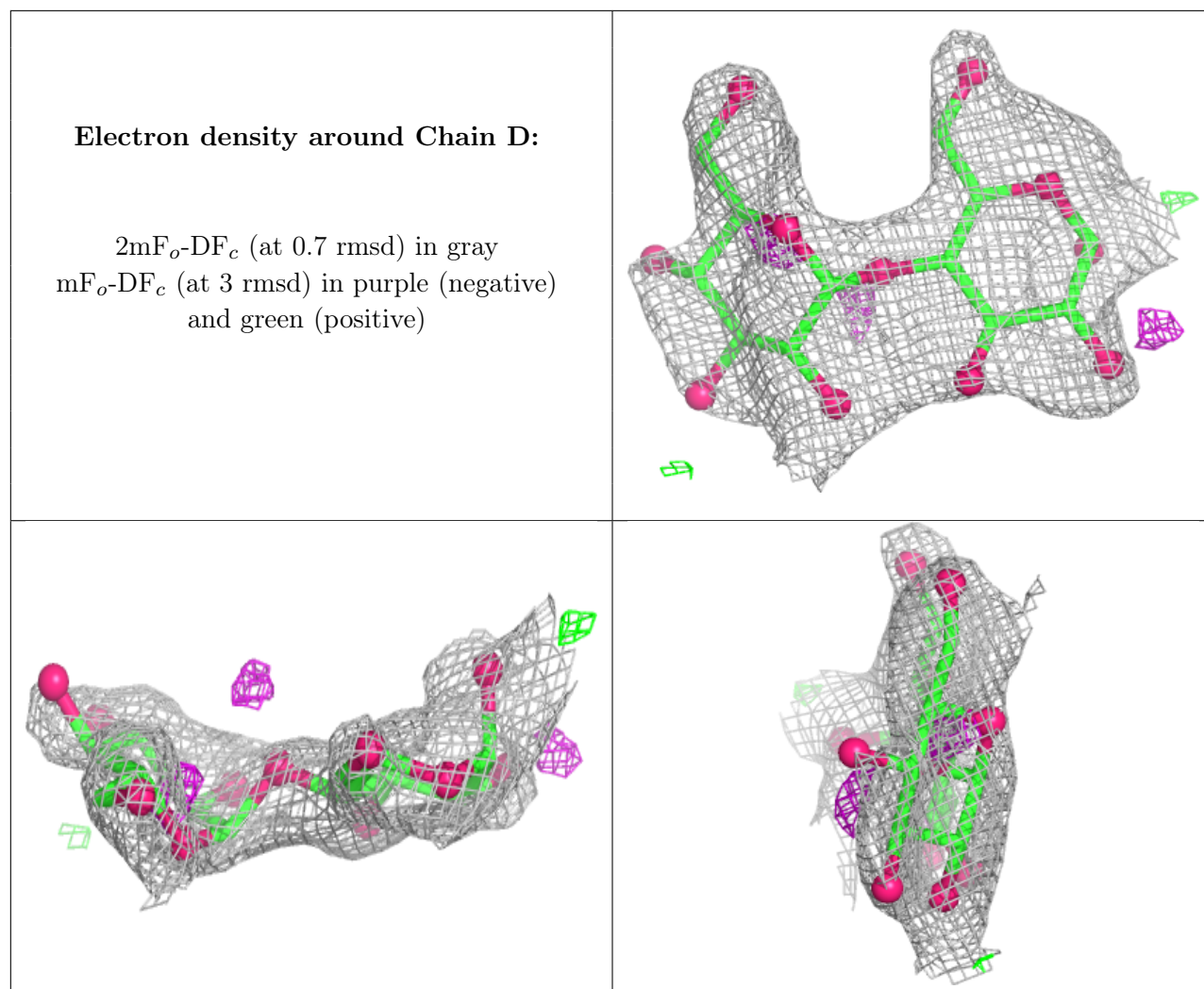
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



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)





6.4 Ligands ⓘ

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
4	CA	A	2405	1/1	0.91	0.16	60,60,60,60	0
4	CA	A	2402	1/1	0.95	0.16	35,35,35,35	0
4	CA	A	2403	1/1	0.98	0.07	30,30,30,30	0
4	CA	A	2404	1/1	0.99	0.04	18,18,18,18	0
4	CA	A	2401	1/1	1.00	0.02	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.