



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:36 PM UTC

PDB ID : 8X8G / pdb_00008x8g
Title : Crystal structure of EndoSz mutant D234M, from Streptococcus equi subsp. Zooepidemicus Sz105, in complex with oligosaccharide G2S2-oxazoline
Authors : Guan, H.H.; Lin, C.C.; Hsieh, Y.C.; Chen, C.J.
Deposited on : 2023-11-27
Resolution : 2.27 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

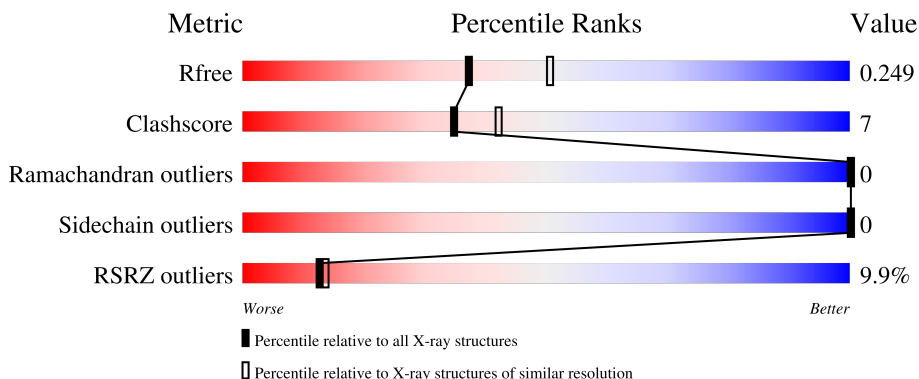
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	961	9% 77% 14% 9%
2	B	9	33% 56% 11%

2 Entry composition [i](#)

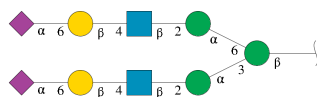
There are 5 unique types of molecules in this entry. The entry contains 7334 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 glycoside hydrolase.

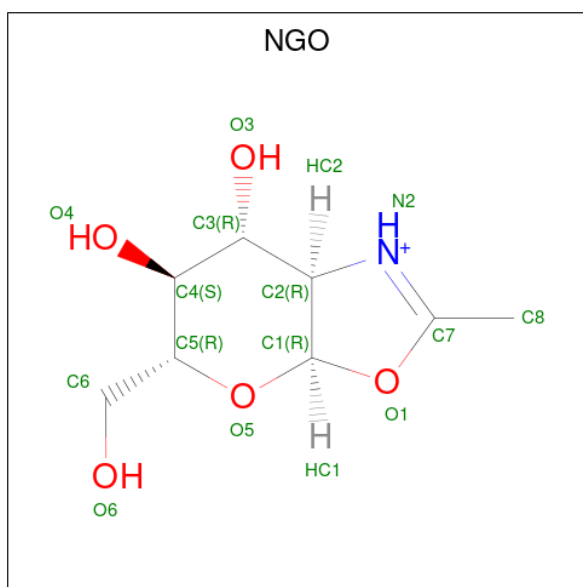
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	876	6902	4375	1152	1360	15	0	0	0

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	9	123	68	4	51	0	0	0

- Molecule 3 is 2-METHYL-4,5-DIHYDRO-(1,2-DIDEOXY-ALPHA-D-GLUCOPYRANOSO)[2,1-D]-1,3-OXAZOLE (CCD ID: NGO) (formula: C₈H₁₄NO₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		

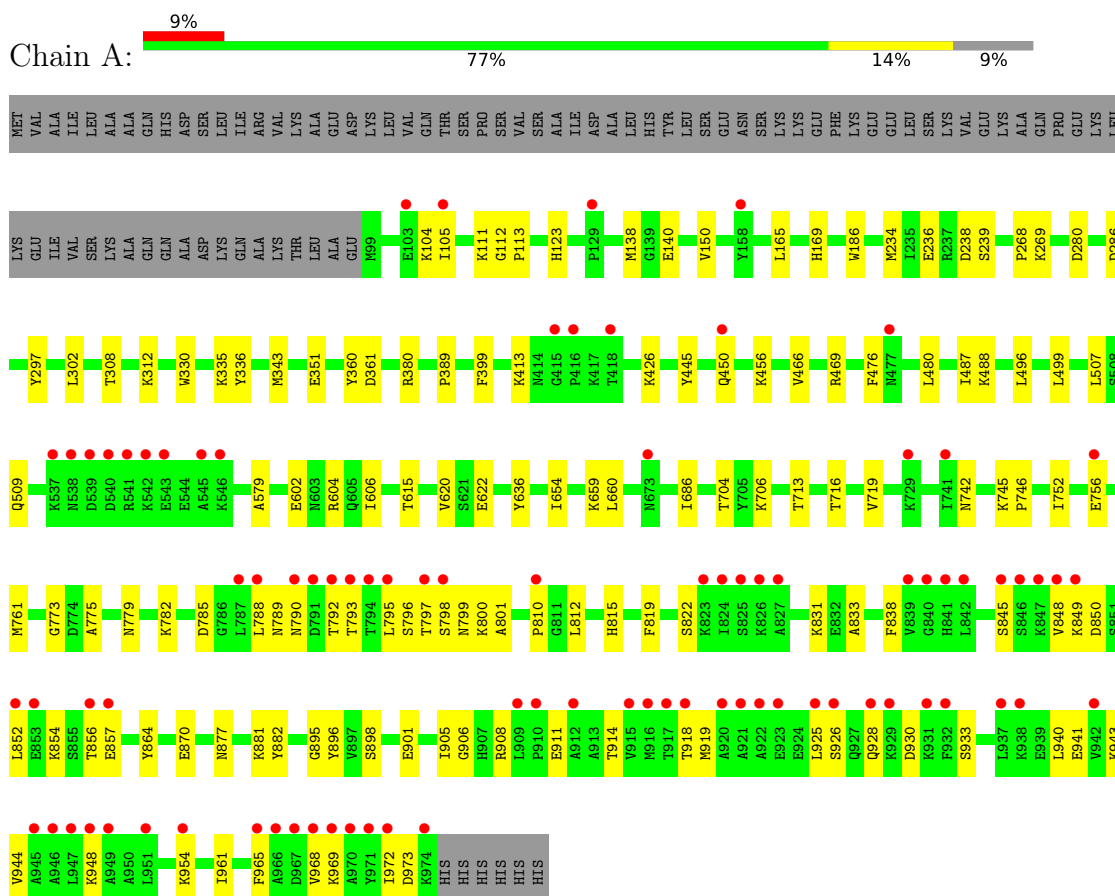
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	294	Total	O	0	0
			294	294		

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: glycoside hydrolase



- Molecule 2: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.06Å 232.60Å 50.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.44 – 2.27 46.44 – 2.27	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.44-2.27) 99.6 (46.44-2.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.27Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.207 , 0.249 0.209 , 0.249	Depositor DCC
R_{free} test set	2846 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7334	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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: NAG, MAN, CA, BMA, GAL, NGO, SIA

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	0.21	0/7035	0.49	0/9501

There are no bond length outliers.

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	6902	0	6887	102	0
2	B	123	0	102	1	0
3	A	14	0	13	2	0
4	A	1	0	0	0	0
5	A	294	0	0	7	0
All	All	7334	0	7002	105	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 7.

All (105) 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:919:MET:HE1	1:A:948:LYS:HD3	1.60	0.83
3:A:10:NGO:HC2	5:A:1126:HOH:O	1.83	0.78
1:A:849:LYS:H	1:A:852:LEU:HD13	1.50	0.75
1:A:941:GLU:HA	1:A:944:VAL:HG12	1.73	0.71
1:A:302:LEU:HD22	1:A:343:MET:HB2	1.76	0.67
1:A:908:ARG:NH2	5:A:1109:HOH:O	2.27	0.66
1:A:848:VAL:O	1:A:849:LYS:HG2	1.96	0.64
3:A:10:NGO:C2	5:A:1126:HOH:O	2.44	0.63
1:A:602:GLU:OE1	1:A:602:GLU:N	2.29	0.62
1:A:713:THR:HG22	1:A:719:VAL:HG22	1.79	0.62
1:A:954:LYS:N	1:A:954:LYS:HD2	2.15	0.61
1:A:286:ASP:N	1:A:286:ASP:OD1	2.32	0.61
1:A:111:LYS:HD2	1:A:112:GLY:N	2.17	0.60
1:A:716:THR:HA	1:A:761:MET:HE3	1.82	0.60
1:A:972:ILE:HG13	1:A:973:ASP:N	2.17	0.59
1:A:782:LYS:NZ	1:A:792:THR:O	2.34	0.59
1:A:450:GLN:H	1:A:450:GLN:CD	2.11	0.59
1:A:789:ASN:OD1	1:A:822:SER:HB2	2.03	0.58
1:A:745:LYS:HD2	1:A:746:PRO:HD2	1.86	0.58
1:A:234:MET:HA	1:A:280:ASP:OD1	2.04	0.57
1:A:831:LYS:NZ	5:A:1117:HOH:O	2.38	0.56
1:A:302:LEU:HB3	1:A:399:PHE:CE1	2.41	0.55
1:A:111:LYS:HE3	1:A:469:ARG:HH21	1.71	0.54
1:A:810:PRO:HG3	1:A:881:LYS:HB3	1.87	0.54
1:A:785:ASP:OD2	1:A:790:ASN:ND2	2.41	0.53
1:A:799:ASN:HB2	1:A:895:GLY:H	1.72	0.53
1:A:919:MET:HE1	1:A:948:LYS:HB2	1.90	0.53
1:A:852:LEU:H	1:A:852:LEU:HD12	1.75	0.52
1:A:833:ALA:HB3	1:A:864:TYR:HB3	1.92	0.52
1:A:845:SER:HB3	1:A:852:LEU:HD11	1.92	0.52
1:A:856:THR:HG22	1:A:857:GLU:N	2.25	0.52
1:A:925:LEU:HA	1:A:928:GLN:HE21	1.75	0.52
1:A:849:LYS:N	1:A:852:LEU:HD13	2.22	0.52
1:A:236:GLU:HG3	1:A:238:ASP:HB3	1.92	0.52
1:A:930:ASP:H	1:A:933:SER:CB	2.22	0.51
1:A:360:TYR:CE1	1:A:413:LYS:HG2	2.45	0.51
1:A:797:THR:OG1	1:A:798:SER:N	2.43	0.51
1:A:795:LEU:HA	1:A:898:SER:HA	1.93	0.50
1:A:961:ILE:O	1:A:965:PHE:N	2.40	0.50
1:A:773:GLY:H	1:A:854:LYS:HE2	1.74	0.50
1:A:845:SER:HB3	1:A:852:LEU:HD21	1.92	0.50
1:A:579:ALA:HB2	1:A:606:ILE:HG23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:838:PHE:HE1	1:A:856:THR:HG21	1.76	0.49
1:A:742:ASN:ND2	5:A:1129:HOH:O	2.45	0.49
1:A:480:LEU:HB3	1:A:499:LEU:HD21	1.94	0.49
1:A:140:GLU:OE2	1:A:426:LYS:NZ	2.44	0.48
1:A:849:LYS:H	1:A:852:LEU:CD1	2.21	0.48
1:A:926:SER:HB2	1:A:940:LEU:HD23	1.95	0.48
1:A:165:LEU:HA	1:A:169:HIS:HB2	1.96	0.48
1:A:812:LEU:HD11	1:A:877:ASN:HB3	1.94	0.48
1:A:280:ASP:HB3	1:A:302:LEU:HB2	1.95	0.48
1:A:756:GLU:OE2	1:A:954:LYS:HG2	2.14	0.47
1:A:269:LYS:NZ	5:A:1119:HOH:O	2.32	0.47
1:A:308:THR:OG1	1:A:312:LYS:HE3	2.15	0.47
1:A:604:ARG:CZ	1:A:686:ILE:HG12	2.44	0.47
1:A:636:TYR:OH	1:A:704:THR:O	2.25	0.47
1:A:819:PHE:CE2	1:A:870:GLU:HG3	2.49	0.47
1:A:848:VAL:C	1:A:849:LYS:HZ2	2.22	0.47
1:A:104:LYS:HD3	1:A:105:ILE:H	1.80	0.46
1:A:972:ILE:HG13	1:A:973:ASP:H	1.80	0.46
1:A:445:TYR:O	1:A:469:ARG:NH1	2.47	0.46
1:A:838:PHE:HB3	1:A:882:TYR:HB2	1.97	0.46
1:A:797:THR:O	1:A:896:TYR:HA	2.16	0.46
1:A:856:THR:HG22	1:A:857:GLU:H	1.81	0.46
1:A:112:GLY:HA2	1:A:113:PRO:C	2.41	0.45
1:A:123:HIS:NE2	1:A:351:GLU:OE2	2.45	0.45
1:A:815:HIS:HB3	1:A:905:ILE:HB	1.98	0.45
1:A:487:ILE:HD12	1:A:507:LEU:HD21	1.99	0.45
1:A:706:LYS:NZ	5:A:1138:HOH:O	2.49	0.45
1:A:186:TRP:HB3	1:A:239:SER:O	2.17	0.44
1:A:330:TRP:CD1	1:A:389:PRO:HB3	2.53	0.44
1:A:775:ALA:HB1	1:A:796:SER:HB3	1.99	0.44
1:A:788:LEU:HD13	1:A:901:GLU:HG3	1.98	0.44
1:A:918:THR:HG21	1:A:965:PHE:HE1	1.82	0.44
1:A:654:ILE:HD12	1:A:659:LYS:HD2	1.99	0.44
1:A:793:THR:HG23	1:A:795:LEU:HD23	1.99	0.44
1:A:797:THR:HG22	1:A:801:ALA:HB2	2.00	0.43
2:B:9:SIA:H6	2:B:9:SIA:H92	1.45	0.43
1:A:234:MET:HE3	1:A:234:MET:HB2	1.64	0.43
1:A:488:LYS:HG2	1:A:509:GLN:OE1	2.19	0.43
1:A:104:LYS:HD3	1:A:104:LYS:HA	1.83	0.43
1:A:111:LYS:HD2	1:A:112:GLY:H	1.81	0.43
1:A:138:MET:HE1	1:A:150:VAL:HG22	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:779:ASN:HA	1:A:782:LYS:HE2	2.00	0.43
1:A:660:LEU:HD11	1:A:752:ILE:HD11	2.00	0.43
1:A:943:LYS:HB3	1:A:968:VAL:HG22	2.00	0.43
1:A:602:GLU:O	1:A:606:ILE:HD12	2.19	0.42
1:A:111:LYS:CE	1:A:469:ARG:HH21	2.33	0.42
1:A:930:ASP:HA	1:A:933:SER:HB2	2.00	0.42
1:A:954:LYS:N	1:A:954:LYS:CD	2.82	0.42
1:A:849:LYS:HG2	1:A:849:LYS:HZ2	1.76	0.42
1:A:911:GLU:OE1	1:A:914:THR:HG23	2.20	0.42
1:A:812:LEU:O	1:A:906:GLY:HA3	2.20	0.41
1:A:852:LEU:HD12	1:A:852:LEU:N	2.36	0.41
1:A:797:THR:HG21	1:A:800:LYS:O	2.20	0.41
1:A:335:LYS:HE2	1:A:336:TYR:CE2	2.56	0.41
1:A:622:GLU:H	1:A:622:GLU:CD	2.27	0.41
1:A:361:ASP:OD1	1:A:380:ARG:HD2	2.20	0.41
1:A:615:THR:HA	1:A:620:VAL:HG21	2.02	0.41
1:A:848:VAL:C	1:A:850:ASP:H	2.28	0.41
1:A:268:PRO:HD2	1:A:297:TYR:O	2.22	0.40
1:A:466:VAL:HG22	1:A:476:PHE:CD2	2.56	0.40
1:A:496:LEU:HD22	1:A:499:LEU:HD13	2.03	0.40
1:A:456:LYS:C	1:A:456:LYS:HD2	2.46	0.40
1:A:969:LYS:O	1:A:972:ILE:HG12	2.21	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	874/961 (91%)	831 (95%)	43 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	753/827 (91%)	753 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	617	HIS
1	A	673	ASN
1	A	763	ASN
1	A	928	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 ⓘ

9 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	BMA	B	1	3,2	11,11,12	0.58	0	15,15,17	0.82	0
2	MAN	B	2	2	11,11,12	0.59	0	15,15,17	1.06	1 (6%)
2	NAG	B	3	2	14,14,15	0.29	0	17,19,21	0.63	0
2	GAL	B	4	2	11,11,12	0.80	0	15,15,17	1.33	2 (13%)
2	SIA	B	5	2	20,20,21	1.65	3 (15%)	21,28,31	1.53	3 (14%)
2	MAN	B	6	2	11,11,12	0.71	0	15,15,17	1.13	2 (13%)
2	NAG	B	7	2	14,14,15	0.40	0	17,19,21	0.59	0
2	GAL	B	8	2	11,11,12	0.59	0	15,15,17	1.09	1 (6%)
2	SIA	B	9	2	20,20,21	1.68	3 (15%)	21,28,31	1.74	6 (28%)

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	BMA	B	1	3,2	-	0/2/19/22	0/1/1/1
2	MAN	B	2	2	-	0/2/19/22	0/1/1/1
2	NAG	B	3	2	-	2/6/23/26	0/1/1/1
2	GAL	B	4	2	-	1/2/19/22	0/1/1/1
2	SIA	B	5	2	-	12/18/34/38	0/1/1/1
2	MAN	B	6	2	-	2/2/19/22	0/1/1/1
2	NAG	B	7	2	-	0/6/23/26	0/1/1/1
2	GAL	B	8	2	-	0/2/19/22	0/1/1/1
2	SIA	B	9	2	-	12/18/34/38	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	5	SIA	C2-C1	5.72	1.59	1.52
2	B	9	SIA	C2-C1	5.70	1.59	1.52
2	B	5	SIA	O6-C2	2.26	1.47	1.43
2	B	9	SIA	C7-C6	2.13	1.55	1.52
2	B	9	SIA	O6-C2	2.13	1.47	1.43
2	B	5	SIA	C7-C6	2.07	1.55	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	9	SIA	C5-N5-C10	5.02	134.86	123.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	5	SIA	O1A-C1-C2	-5.01	112.03	122.85
2	B	4	GAL	C1-O5-C5	3.53	116.92	112.19
2	B	2	MAN	C1-O5-C5	2.96	116.15	112.19
2	B	6	MAN	O2-C2-C3	-2.86	104.22	110.15
2	B	9	SIA	C4-C5-N5	2.59	115.54	110.44
2	B	6	MAN	C1-O5-C5	2.59	115.66	112.19
2	B	9	SIA	O1A-C1-C2	-2.49	117.46	122.85
2	B	8	GAL	C1-O5-C5	2.48	115.51	112.19
2	B	9	SIA	O1B-C1-O1A	2.37	129.46	124.08
2	B	9	SIA	C11-C10-N5	2.34	120.00	116.12
2	B	5	SIA	O1B-C1-O1A	2.33	129.37	124.08
2	B	5	SIA	O1B-C1-C2	2.27	118.61	112.71
2	B	4	GAL	O2-C2-C3	-2.10	105.81	110.15
2	B	9	SIA	O10-C10-C11	-2.09	118.33	122.05

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	5	SIA	C5-C6-C7-C8
2	B	5	SIA	C5-C6-C7-O7
2	B	5	SIA	O6-C6-C7-C8
2	B	5	SIA	O6-C6-C7-O7
2	B	5	SIA	O8-C8-C9-O9
2	B	9	SIA	C6-C7-C8-C9
2	B	9	SIA	C6-C7-C8-O8
2	B	9	SIA	O7-C7-C8-O8
2	B	9	SIA	C7-C8-C9-O9
2	B	9	SIA	O8-C8-C9-O9
2	B	3	NAG	C4-C5-C6-O6
2	B	5	SIA	C7-C8-C9-O9
2	B	3	NAG	O5-C5-C6-O6
2	B	6	MAN	C4-C5-C6-O6
2	B	6	MAN	O5-C5-C6-O6
2	B	5	SIA	C11-C10-N5-C5
2	B	5	SIA	O10-C10-N5-C5
2	B	9	SIA	C11-C10-N5-C5
2	B	9	SIA	O10-C10-N5-C5
2	B	9	SIA	O7-C7-C8-C9
2	B	4	GAL	O5-C5-C6-O6
2	B	5	SIA	C6-C7-C8-O8
2	B	9	SIA	C6-C5-N5-C10

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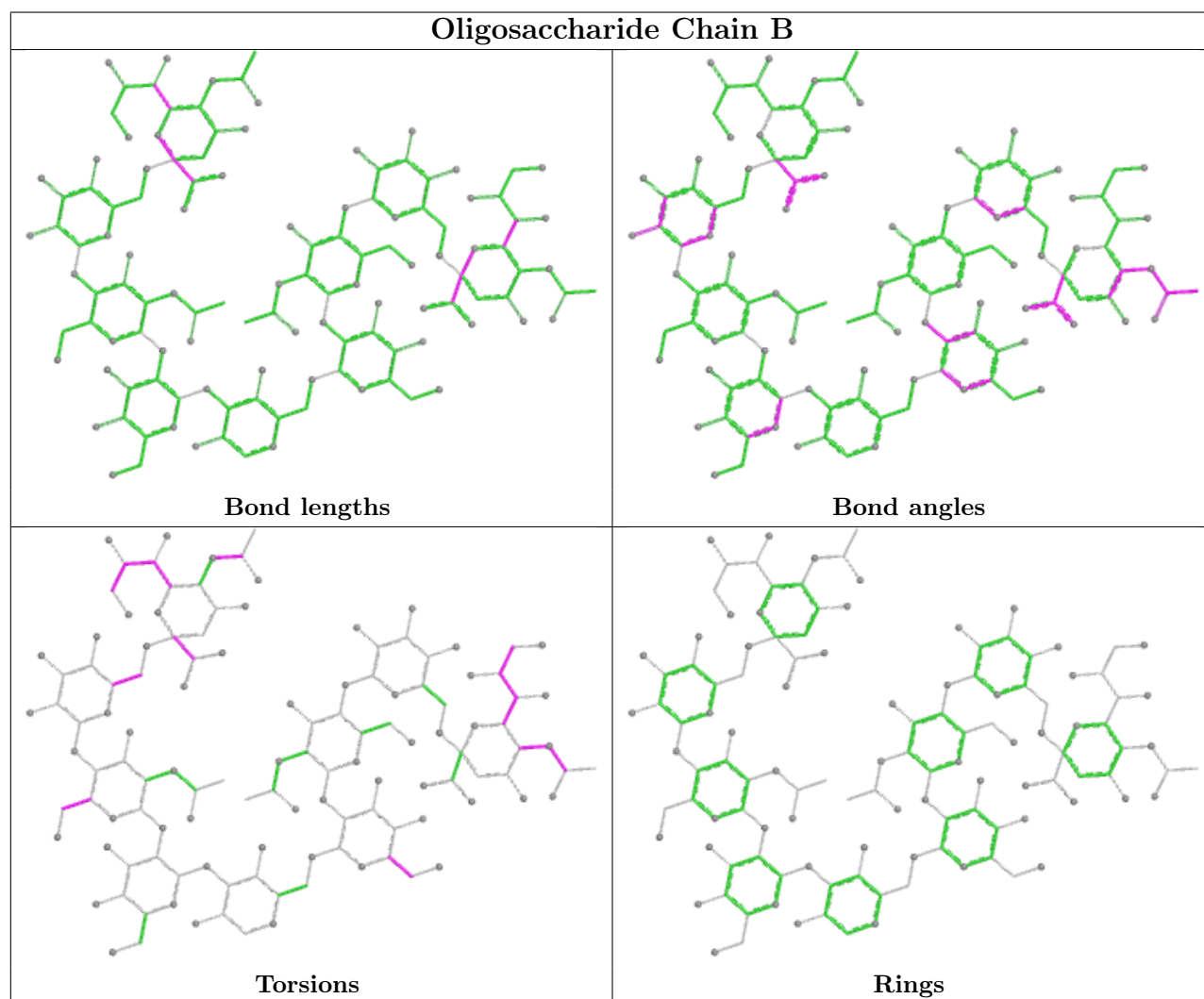
Mol	Chain	Res	Type	Atoms
2	B	9	SIA	O6-C6-C7-O7
2	B	5	SIA	C6-C7-C8-C9
2	B	9	SIA	C5-C6-C7-C8
2	B	5	SIA	O1A-C1-C2-O6
2	B	5	SIA	O7-C7-C8-O8
2	B	9	SIA	C4-C5-N5-C10

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	9	SIA	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

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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
3	NGO	A	10	2	14,15,15	3.61	6 (42%)	15,22,22	2.90	7 (46%)

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
3	NGO	A	10	2	-	1/2/30/30	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	10	NGO	C1-C2	-8.51	1.36	1.52
3	A	10	NGO	O5-C1	7.73	1.61	1.41
3	A	10	NGO	C2-N2	4.64	1.52	1.47
3	A	10	NGO	C4-C5	-2.18	1.48	1.53
3	A	10	NGO	O5-C5	-2.15	1.39	1.44
3	A	10	NGO	C3-C2	2.11	1.57	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	10	NGO	C3-C2-N2	-5.19	103.29	110.56
3	A	10	NGO	C1-O1-C7	-4.73	100.35	106.88
3	A	10	NGO	C1-O5-C5	-4.54	104.86	113.72
3	A	10	NGO	O5-C1-O1	4.42	117.01	109.25
3	A	10	NGO	C2-N2-C7	-3.84	103.75	106.43
3	A	10	NGO	O6-C6-C5	-3.38	99.82	111.33
3	A	10	NGO	C8-C7-N2	-2.25	123.81	127.05

There are no chirality outliers.

All (1) torsion outliers are listed below:

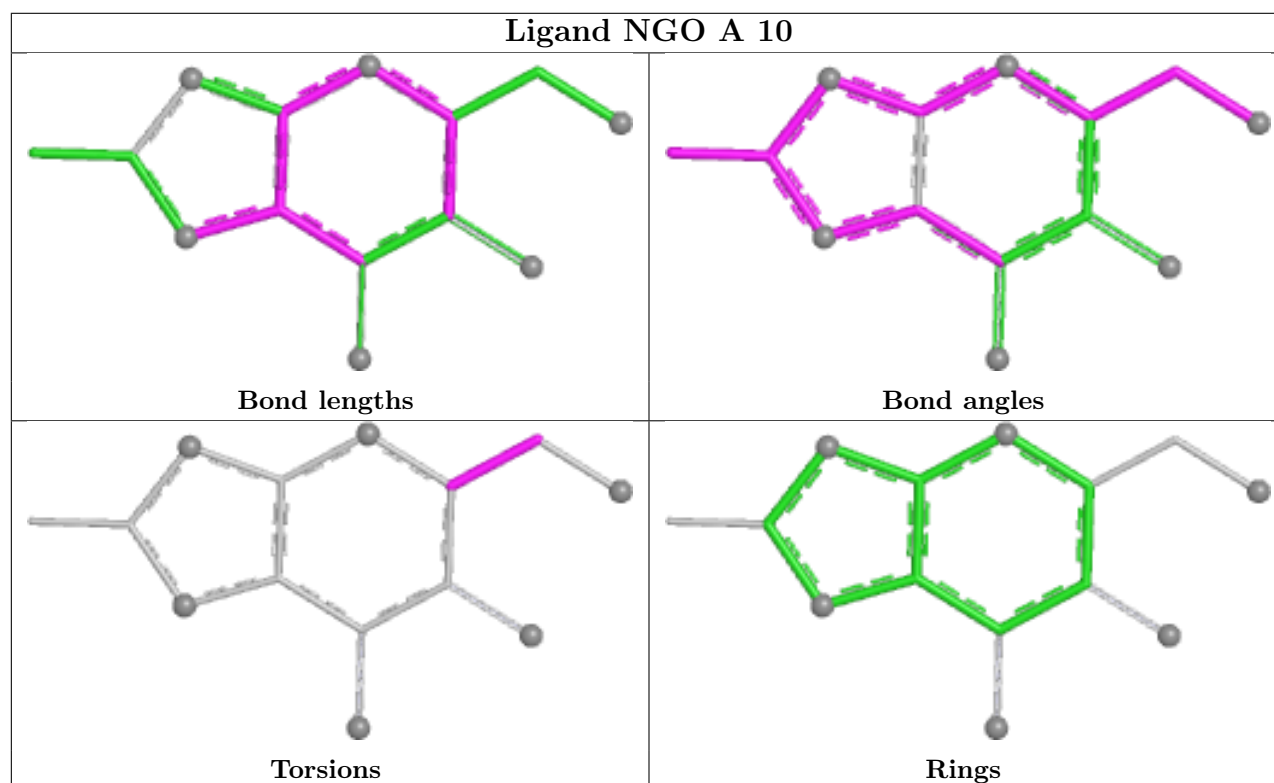
Mol	Chain	Res	Type	Atoms
3	A	10	NGO	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	10	NGO	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 ⓘ

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	876/961 (91%)	0.55	87 (9%) 13 14	28, 51, 108, 161	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	798	SER	6.0
1	A	842	LEU	5.9
1	A	793	THR	4.7
1	A	792	THR	4.2
1	A	972	ILE	4.1
1	A	839	VAL	4.0
1	A	788	LEU	3.9
1	A	826	LYS	3.8
1	A	848	VAL	3.6
1	A	824	ILE	3.6
1	A	937	LEU	3.6
1	A	932	PHE	3.5
1	A	912	ALA	3.5
1	A	925	LEU	3.5
1	A	942	VAL	3.4
1	A	846	SER	3.4
1	A	827	ALA	3.4
1	A	968	VAL	3.4
1	A	537	LYS	3.3
1	A	856	THR	3.3
1	A	852	LEU	3.2
1	A	910	PRO	3.1
1	A	931	LYS	3.0
1	A	915	VAL	3.0
1	A	105	ILE	3.0
1	A	416	PRO	3.0
1	A	477	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	909	LEU	2.9
1	A	840	GLY	2.9
1	A	841	HIS	2.9
1	A	540	ASP	2.9
1	A	810	PRO	2.8
1	A	539	ASP	2.8
1	A	966	ALA	2.8
1	A	970	ALA	2.8
1	A	825	SER	2.8
1	A	922	ALA	2.8
1	A	945	ALA	2.8
1	A	965	PHE	2.7
1	A	790	ASN	2.7
1	A	847	LYS	2.7
1	A	541	ARG	2.7
1	A	853	GLU	2.7
1	A	158	TYR	2.7
1	A	545	ALA	2.6
1	A	946	ALA	2.6
1	A	857	GLU	2.6
1	A	938	LYS	2.6
1	A	794	THR	2.6
1	A	918	THR	2.6
1	A	795	LEU	2.6
1	A	791	ASP	2.6
1	A	974	LYS	2.5
1	A	756	GLU	2.5
1	A	538	ASN	2.5
1	A	929	LYS	2.5
1	A	967	ASP	2.4
1	A	951	LEU	2.4
1	A	928	GLN	2.4
1	A	921	ALA	2.4
1	A	103	GLU	2.4
1	A	845	SER	2.3
1	A	741	ILE	2.3
1	A	673	ASN	2.3
1	A	849	LYS	2.3
1	A	415	GLY	2.3
1	A	916	MET	2.3
1	A	418	THR	2.3
1	A	948	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	542	LYS	2.2
1	A	969	LYS	2.2
1	A	450	GLN	2.2
1	A	823	LYS	2.2
1	A	920	ALA	2.2
1	A	917	THR	2.2
1	A	129	PRO	2.2
1	A	787	LEU	2.1
1	A	546	LYS	2.1
1	A	797	THR	2.1
1	A	926	SER	2.1
1	A	949	ALA	2.1
1	A	923	GLU	2.1
1	A	954	LYS	2.0
1	A	947	LEU	2.0
1	A	971	TYR	2.0
1	A	729	LYS	2.0
1	A	543	GLU	2.0

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

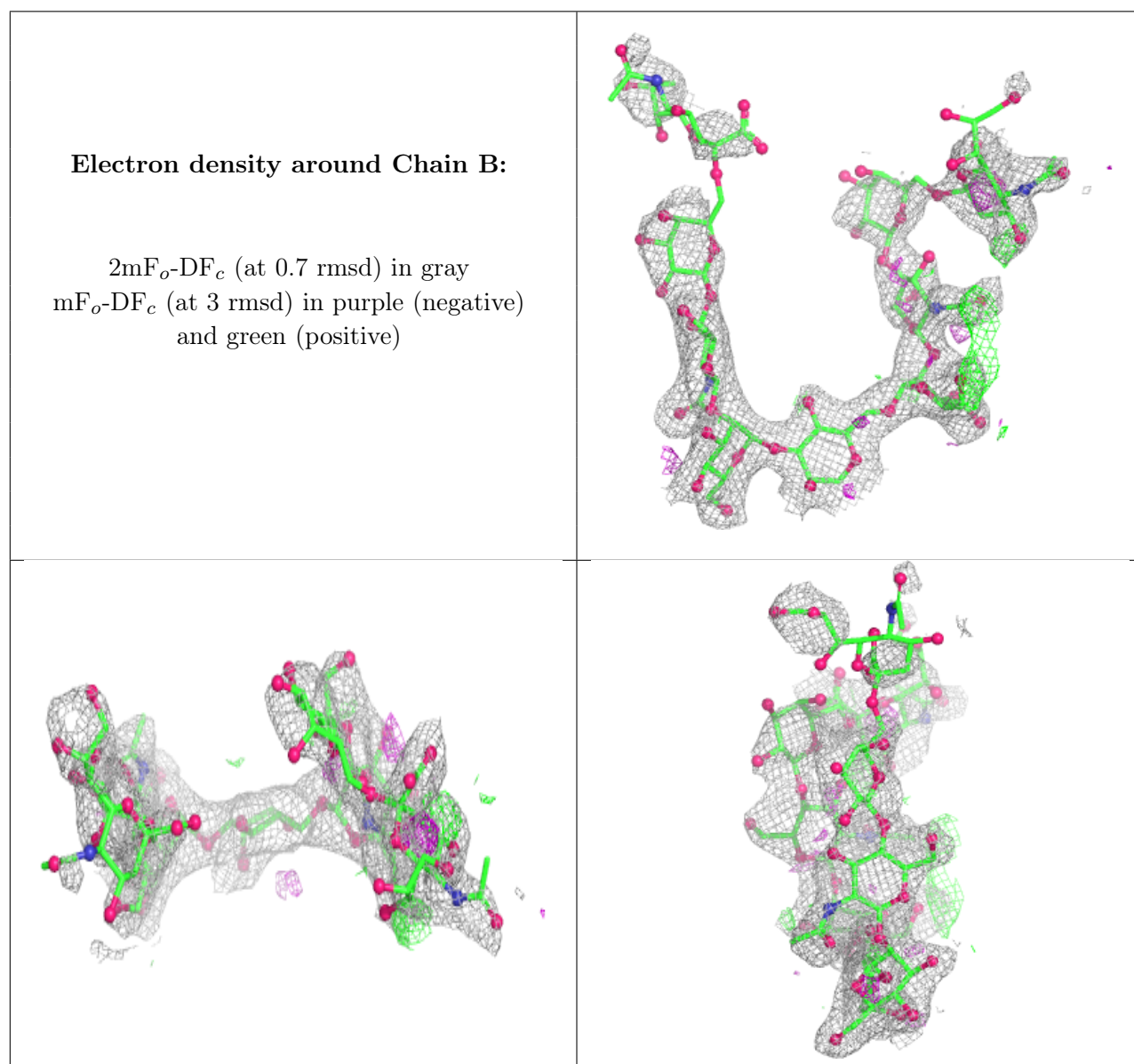
There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
2	SIA	B	5	20/21	0.56	0.21	88,129,138,138	0
2	GAL	B	4	11/12	0.58	0.17	80,99,119,131	0
2	GAL	B	8	11/12	0.58	0.16	109,114,118,119	0
2	SIA	B	9	20/21	0.65	0.20	92,106,118,118	0
2	NAG	B	7	14/15	0.67	0.22	72,89,102,111	0
2	MAN	B	6	11/12	0.79	0.15	55,63,67,71	0
2	NAG	B	3	14/15	0.91	0.11	48,62,69,80	0
2	MAN	B	2	11/12	0.96	0.07	36,39,48,51	0
2	BMA	B	1	11/12	0.96	0.07	31,40,53,55	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 ⓘ

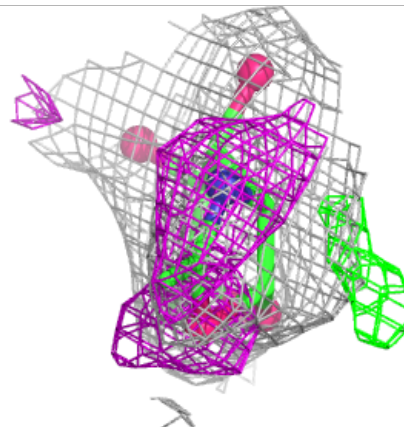
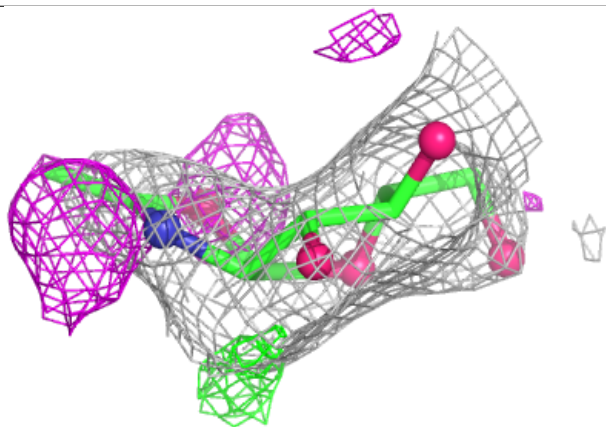
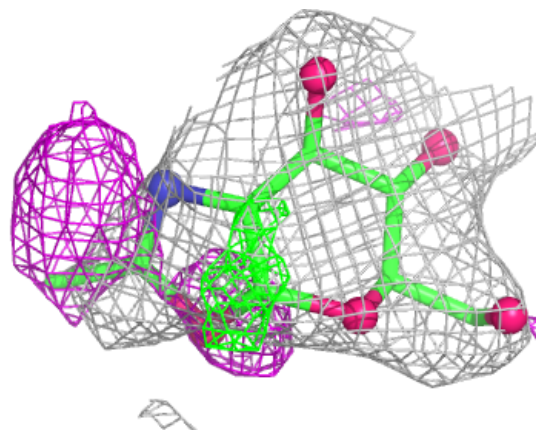
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NGO	A	10	14/14	0.88	0.13	39,55,65,66	0
4	CA	A	1002	1/1	0.91	0.12	82,82,82,82	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

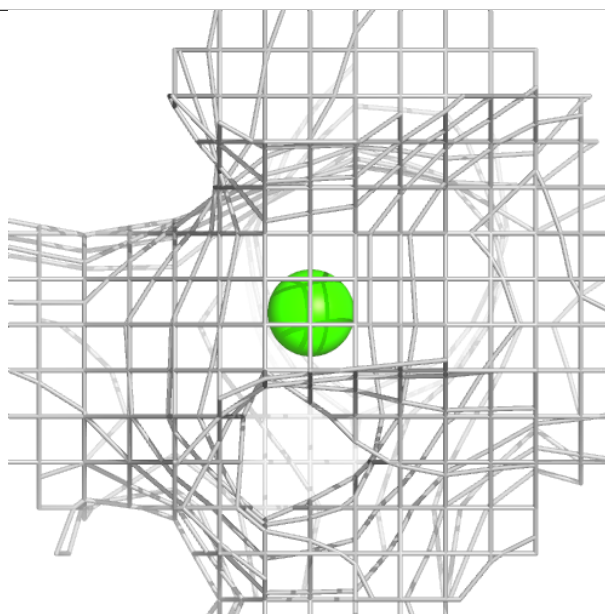
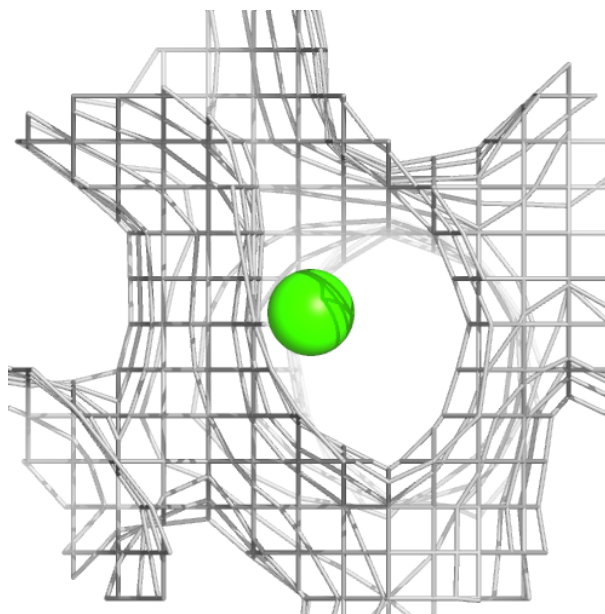
Electron density around NGO A 10:

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



Electron density around CA A 1002:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.