



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:23 PM UTC

PDB ID : 4ZPP / pdb_00004zpp
Title : Crystal Structure of Protocadherin Gamma C5 EC1-3
Authors : Wolcott, H.N.; Goodman, K.M.; Bahna, F.; Mannepalli, S.; Shapiro, L.
Deposited on : 2015-05-08
Resolution : 3.00 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

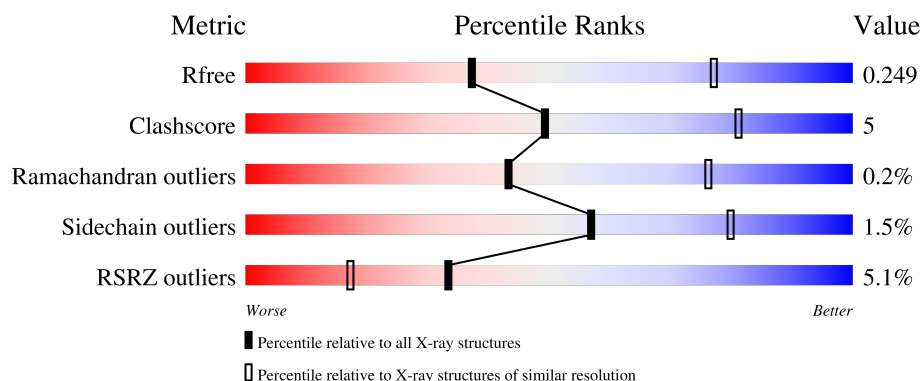
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	324	2% 86% 10% .
1	B	324	8% 84% 12% .
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4614 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 MCG133388, isoform CRA_f.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	3	0
			2327	1437	408	477	5			
1	B	311	Total	C	N	O	S	0	0	0
			2132	1296	377	454	5			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	317	HIS	-	expression tag	UNP Q91XW9
A	318	HIS	-	expression tag	UNP Q91XW9
A	319	HIS	-	expression tag	UNP Q91XW9
A	320	HIS	-	expression tag	UNP Q91XW9
A	321	HIS	-	expression tag	UNP Q91XW9
A	322	HIS	-	expression tag	UNP Q91XW9
A	323	HIS	-	expression tag	UNP Q91XW9
A	324	HIS	-	expression tag	UNP Q91XW9
B	317	HIS	-	expression tag	UNP Q91XW9
B	318	HIS	-	expression tag	UNP Q91XW9
B	319	HIS	-	expression tag	UNP Q91XW9
B	320	HIS	-	expression tag	UNP Q91XW9
B	321	HIS	-	expression tag	UNP Q91XW9
B	322	HIS	-	expression tag	UNP Q91XW9
B	323	HIS	-	expression tag	UNP Q91XW9
B	324	HIS	-	expression tag	UNP Q91XW9

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

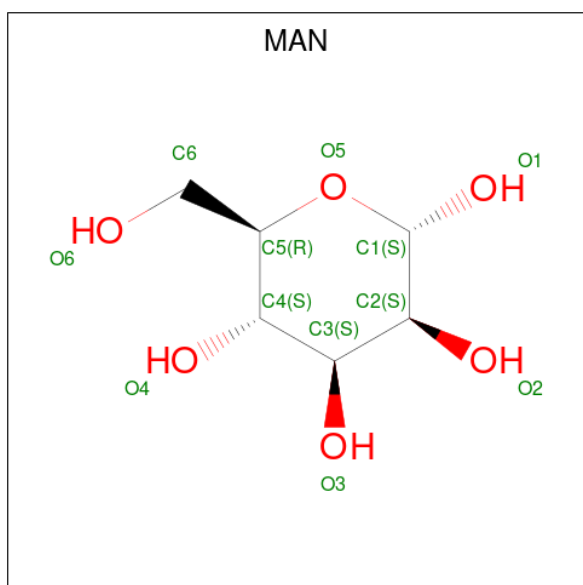


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	Ca	0	0
			6	6		
3	B	6	Total	Ca	0	0
			6	6		

- Molecule 4 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	A	1	Total	C	O	0	0
			11	6	5		
4	B	1	Total	C	O	0	0
			11	6	5		
4	B	1	Total	C	O	0	0
			11	6	5		

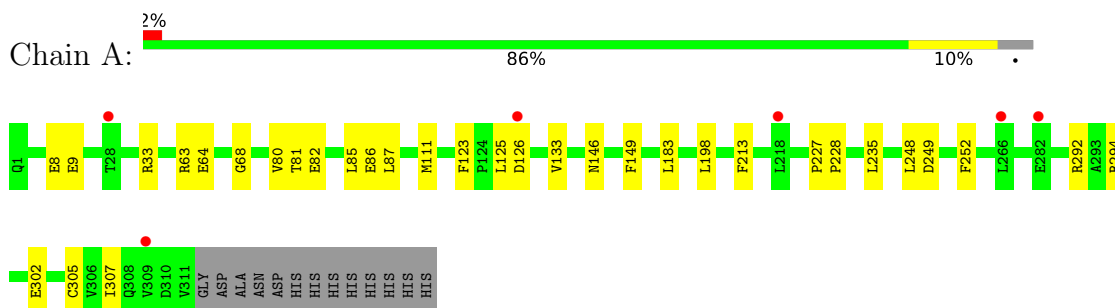
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	34	Total 34	O 34	0	0
5	B	9	Total 9	O 9	0	0

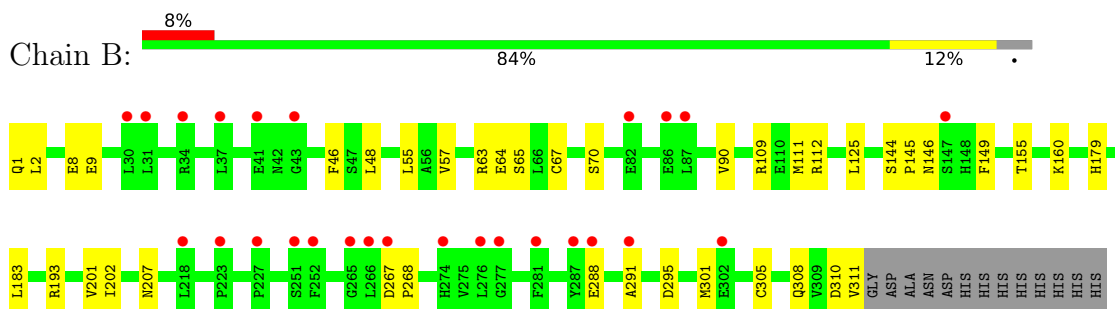
3 Residue-property plots [i](#)

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: MCG133388, isoform CRA_f



- Molecule 1: MCG133388, isoform CRA_f



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.19Å 84.56Å 109.14Å 90.00° 106.43° 90.00°	Depositor
Resolution (Å)	27.94 – 3.00 27.94 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (27.94-3.00) 94.8 (27.94-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 3.00Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.209 , 0.248 0.220 , 0.249	Depositor DCC
R_{free} test set	1202 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	69.7	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 74.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4614	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% 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, NAG, MAN

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	0.28	0/2377	0.71	0/3243
1	B	0.28	0/2168	0.73	0/2974
All	All	0.28	0/4545	0.72	0/6217

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

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	2327	0	2215	23	0
1	B	2132	0	1825	24	0
2	C	28	0	25	1	0
2	D	28	0	25	1	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
4	A	22	0	20	1	0
4	B	22	0	20	0	0
5	A	34	0	0	0	0
5	B	9	0	0	0	0
All	All	4614	0	4130	47	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 5.

All (47) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:HH12	1:B:111:MET:HE3	1.59	0.68
1:A:235:LEU:HD21	1:A:307:ILE:HD11	1.78	0.66
1:B:8:GLU:HG2	1:B:9:GLU:HG2	1.81	0.63
1:A:9:GLU:OE2	1:B:193:ARG:NH2	2.35	0.60
1:A:8:GLU:HG2	1:A:9:GLU:HG2	1.84	0.59
1:B:179:HIS:HB2	1:B:201:VAL:HB	1.83	0.59
1:B:8:GLU:HG3	1:B:63:ARG:H	1.69	0.58
1:B:146:ASN:HD21	1:B:149:PHE:HB2	1.69	0.57
2:D:1:NAG:H62	2:D:2:NAG:H83	1.85	0.57
1:A:63:ARG:NH1	1:A:64:GLU:OE1	2.39	0.55
1:A:63:ARG:NH2	1:A:68:GLY:O	2.36	0.55
2:C:1:NAG:H4	2:C:2:NAG:N2	2.22	0.55
1:A:8:GLU:HG3	1:A:63:ARG:H	1.70	0.55
1:B:46:PHE:HB3	1:B:55:LEU:HD11	1.89	0.54
1:A:111:MET:HE1	1:A:123:PHE:HB3	1.89	0.54
1:B:295:ASP:HB3	1:B:301:MET:H	1.73	0.53
1:A:146:ASN:HD21	1:A:149:PHE:HB2	1.72	0.53
1:B:288:GLU:HG2	1:B:308:GLN:HG2	1.93	0.51
1:B:63:ARG:NH1	1:B:64:GLU:OE1	2.46	0.49
1:B:125:LEU:HD13	1:B:183:LEU:HD22	1.93	0.49
1:B:112:ARG:HG2	1:B:202:ILE:HB	1.93	0.49
1:B:291:ALA:HB3	1:B:305:CYS:HB3	1.95	0.49
1:A:80:VAL:HG22	1:A:87:LEU:HD22	1.96	0.47
1:A:125:LEU:HD13	1:A:183:LEU:HD22	1.95	0.47
1:B:146:ASN:ND2	1:B:149:PHE:H	2.12	0.47
1:A:126:ASP:OD1	1:A:126:ASP:N	2.46	0.47
1:A:213:PHE:CD1	1:A:305:CYS:HB2	2.50	0.47
1:A:85:LEU:H	1:A:85:LEU:HD23	1.81	0.46
4:A:409:MAN:H3	1:B:65:SER:HB3	1.99	0.45
1:A:33[B]:ARG:NH2	1:A:86:GLU:OE1	2.37	0.45
1:A:146:ASN:ND2	1:A:149:PHE:H	2.15	0.45
1:A:252:PHE:O	1:A:292:ARG:NH2	2.50	0.45
1:B:2:LEU:HB2	1:B:90:VAL:HG22	1.98	0.44
1:B:8:GLU:HG2	1:B:9:GLU:N	2.32	0.44
1:B:146:ASN:HD21	1:B:149:PHE:H	1.66	0.43
1:A:9:GLU:CD	1:B:193:ARG:HH22	2.26	0.43
1:A:294:ARG:NH1	1:A:302:GLU:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:MET:HE2	1:A:111:MET:HB3	1.94	0.43
1:B:267:ASP:HA	1:B:268:PRO:HD3	1.90	0.42
1:B:310:ASP:OD1	1:B:311:VAL:N	2.46	0.41
1:A:81:THR:O	1:A:85:LEU:HA	2.20	0.41
1:A:227:PRO:HA	1:A:228:PRO:HD3	1.92	0.41
1:B:144:SER:HA	1:B:145:PRO:HD3	1.97	0.41
1:A:249:ASP:N	1:A:249:ASP:OD1	2.54	0.41
1:B:160:LYS:HA	1:B:160:LYS:HD2	1.86	0.41
1:A:82:GLU:HA	1:A:85:LEU:HA	2.03	0.41
1:B:67:CYS:HB3	1:B:70:SER:HB2	2.02	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	312/324 (96%)	304 (97%)	8 (3%)	0	100	100
1	B	309/324 (95%)	291 (94%)	17 (6%)	1 (0%)	36	70
All	All	621/648 (96%)	595 (96%)	25 (4%)	1 (0%)	43	76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	207	ASN

5.3.2 Protein sidechains [i](#)

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	256/280 (91%)	253 (99%)	3 (1%)	63	82
1	B	205/280 (73%)	201 (98%)	4 (2%)	48	76
All	All	461/560 (82%)	454 (98%)	7 (2%)	57	80

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	VAL
1	A	198	LEU
1	A	248	LEU
1	B	1	GLN
1	B	48	LEU
1	B	57	VAL
1	B	155	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	ASN
1	A	152	HIS
1	A	209	ASN
1	A	274	HIS
1	B	146	ASN
1	B	209	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	NAG	C	1	2,1	14,14,15	0.25	0	17,19,21	0.43	0
2	NAG	C	2	2	14,14,15	0.61	1 (7%)	17,19,21	0.65	1 (5%)
2	NAG	D	1	2,1	14,14,15	0.66	1 (7%)	17,19,21	0.97	1 (5%)
2	NAG	D	2	2	14,14,15	0.37	0	17,19,21	0.60	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	NAG	C	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	1/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NAG	O5-C1	2.17	1.47	1.43
2	C	2	NAG	C1-C2	2.09	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	NAG	C1-O5-C5	3.74	117.20	112.19
2	C	2	NAG	C1-O5-C5	2.13	115.03	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	O5-C5-C6-O6

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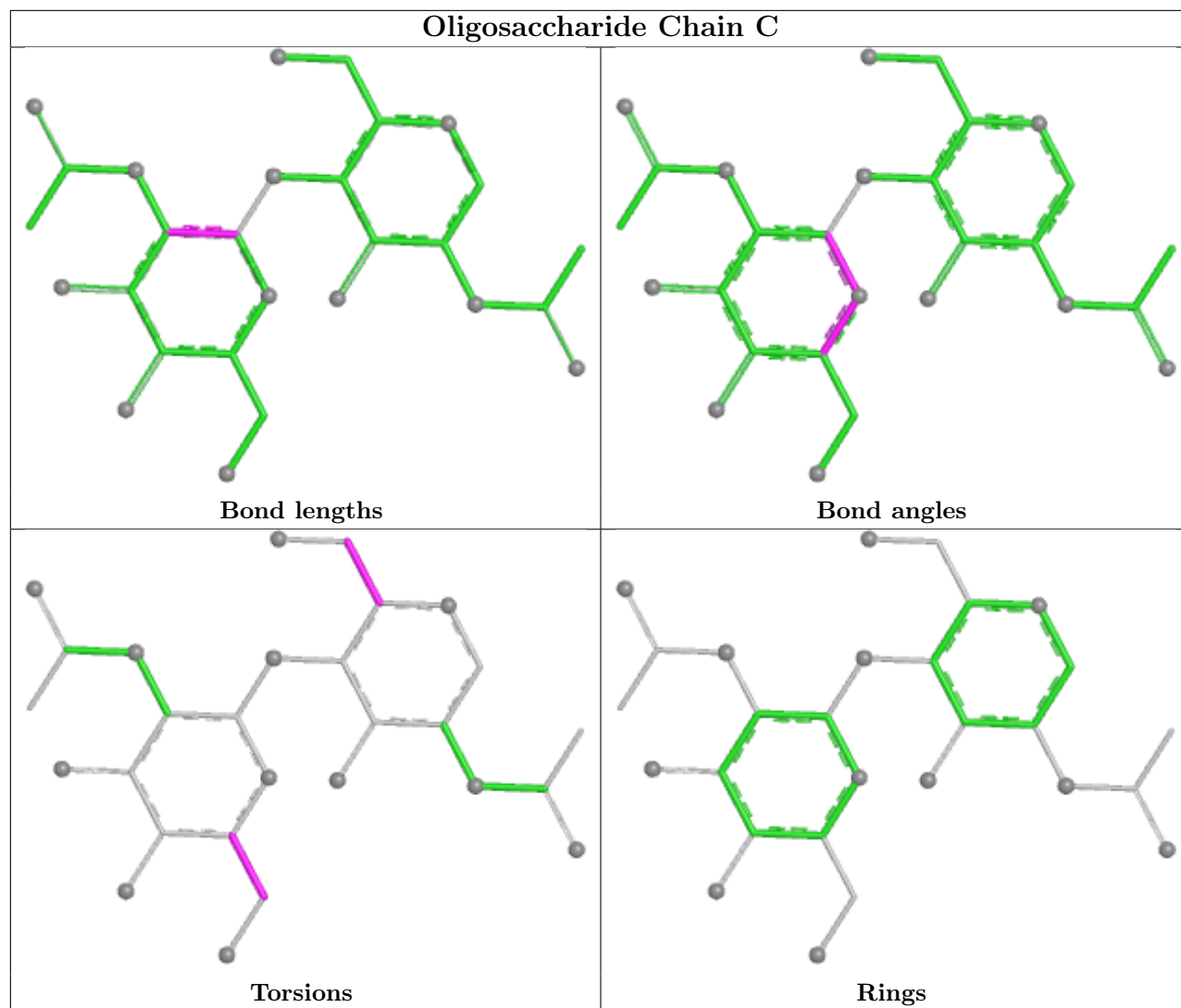
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C3-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7

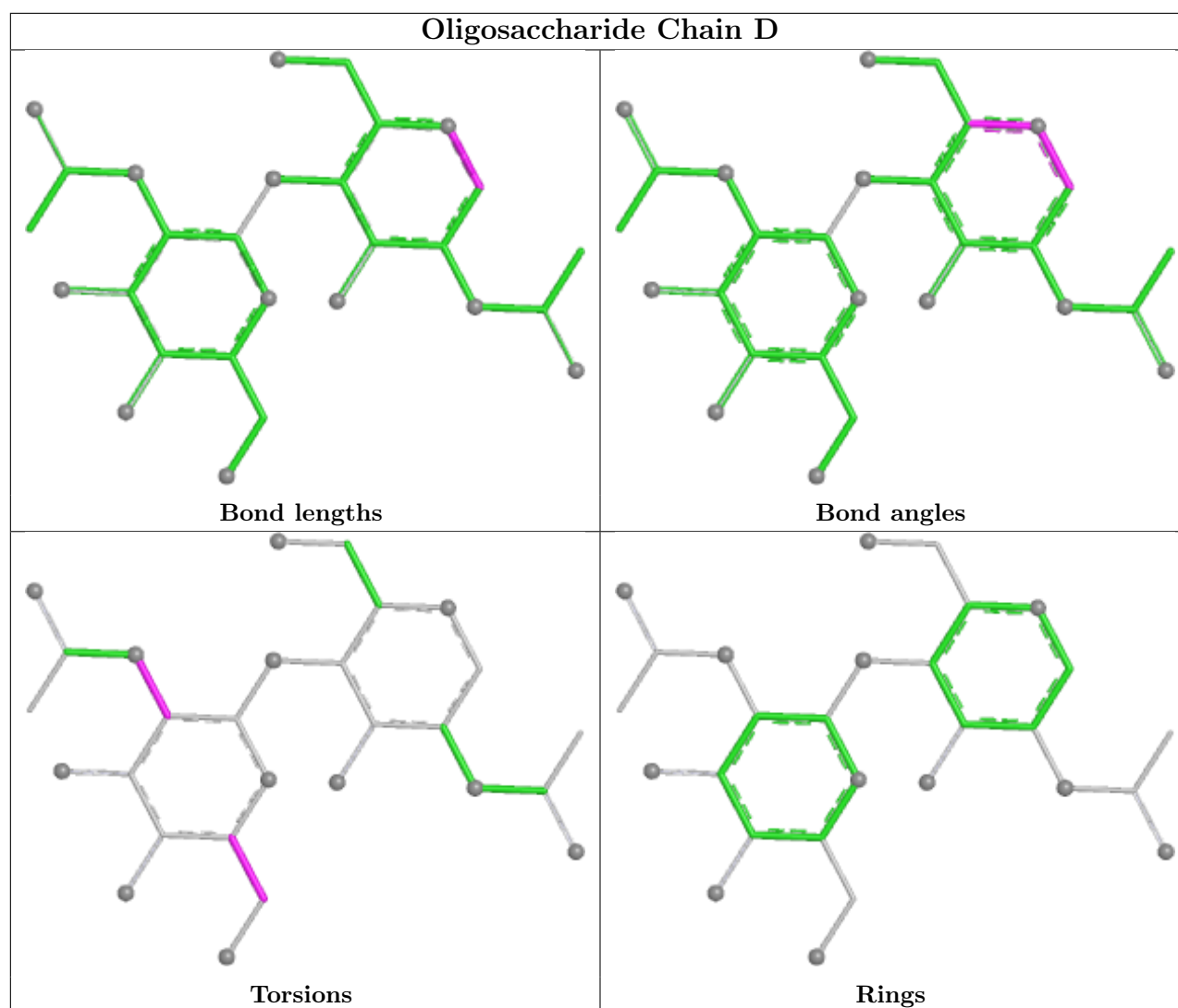
There are no ring outliers.

4 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	C	2	NAG	1	0
2	D	2	NAG	1	0
2	D	1	NAG	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 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	MAN	A	409	1	11,11,12	0.61	0	15,15,17	1.13	2 (13%)
4	MAN	A	410	1	11,11,12	0.61	0	15,15,17	1.01	2 (13%)
4	MAN	B	410	1	11,11,12	0.87	1 (9%)	15,15,17	0.87	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	B	409	1	11,11,12	0.54	0	15,15,17	1.15	2 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	A	409	1	-	2/2/19/22	0/1/1/1
4	MAN	A	410	1	-	0/2/19/22	0/1/1/1
4	MAN	B	410	1	-	2/2/19/22	0/1/1/1
4	MAN	B	409	1	-	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)
4	B	410	MAN	O5-C1	-2.26	1.39	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	409	MAN	C1-O5-C5	3.26	116.56	112.19
4	B	409	MAN	C1-O5-C5	3.03	116.25	112.19
4	A	410	MAN	C1-O5-C5	2.49	115.53	112.19
4	B	409	MAN	O2-C2-C3	-2.35	105.29	110.15
4	A	410	MAN	O2-C2-C3	-2.20	105.60	110.15
4	B	410	MAN	O2-C2-C3	-2.12	105.77	110.15
4	A	409	MAN	O2-C2-C3	-2.03	105.94	110.15

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	410	MAN	O5-C5-C6-O6
4	B	410	MAN	C4-C5-C6-O6
4	A	409	MAN	O5-C5-C6-O6
4	A	409	MAN	C4-C5-C6-O6
4	B	409	MAN	O5-C5-C6-O6
4	B	409	MAN	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	409	MAN	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	311/324 (95%)	0.04	6 (1%) 66 43	37, 81, 166, 251	3 (0%)
1	B	311/324 (95%)	0.43	26 (8%) 17 9	43, 121, 211, 263	0
All	All	622/648 (95%)	0.24	32 (5%) 33 17	37, 92, 198, 263	3 (0%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	282	GLU	5.3
1	B	252	PHE	3.8
1	A	266	LEU	3.8
1	A	126	ASP	3.8
1	B	266	LEU	3.5
1	B	43	GLY	3.5
1	B	281	PHE	3.4
1	B	227	PRO	3.4
1	B	41	GLU	3.2
1	B	87	LEU	3.1
1	B	30	LEU	3.1
1	B	274	HIS	3.0
1	B	218	LEU	3.0
1	B	287	TYR	3.0
1	B	86	GLU	2.7
1	B	265	GLY	2.6
1	B	302	GLU	2.6
1	B	288	GLU	2.5
1	B	223	PRO	2.5
1	B	291	ALA	2.4
1	A	309	VAL	2.4
1	B	82	GLU	2.4
1	B	251	SER	2.3
1	B	31	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	218	LEU	2.2
1	A	28	THR	2.1
1	B	276	LEU	2.1
1	B	34	ARG	2.1
1	B	37	LEU	2.1
1	B	267	ASP	2.1
1	B	277	GLY	2.1
1	B	147	SER	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](#)

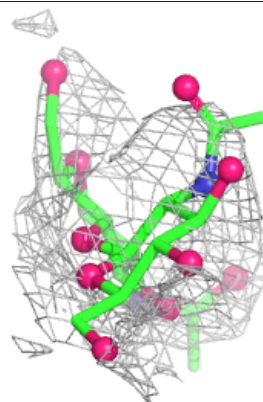
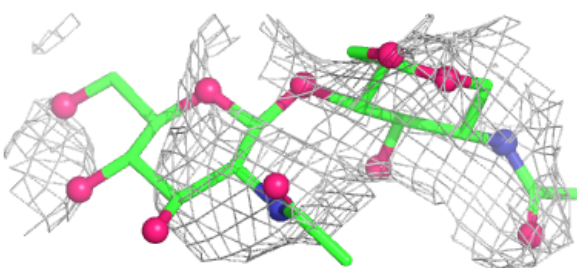
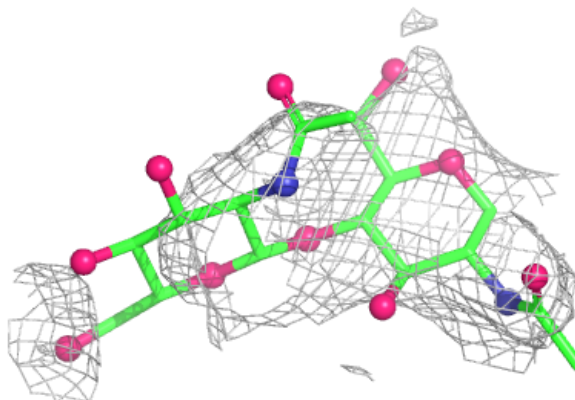
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NAG	C	2	14/15	0.32	0.16	159,180,186,186	0
2	NAG	D	2	14/15	0.58	0.13	170,180,184,184	0
2	NAG	C	1	14/15	0.66	0.15	121,138,158,169	0
2	NAG	D	1	14/15	0.70	0.13	123,152,161,171	0

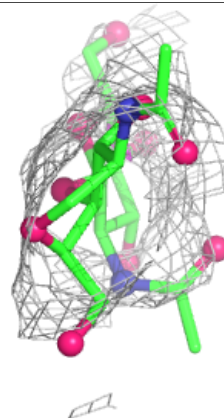
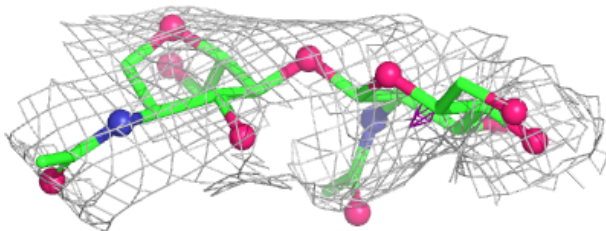
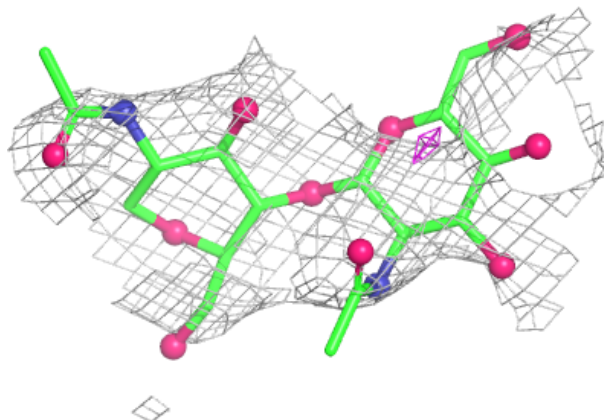
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

**Electron density around Chain D:**

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



6.4 Ligands

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	MAN	A	409	11/12	0.89	0.16	80,104,115,121	0
4	MAN	B	409	11/12	0.91	0.15	70,80,102,106	0
4	MAN	B	410	11/12	0.92	0.11	62,70,112,113	0
4	MAN	A	410	11/12	0.93	0.09	78,97,110,114	0
3	CA	B	402	1/1	0.96	0.08	88,88,88,88	0
3	CA	A	406	1/1	0.98	0.09	63,63,63,63	0
3	CA	B	405	1/1	0.99	0.04	57,57,57,57	0
3	CA	B	406	1/1	0.99	0.05	57,57,57,57	0
3	CA	A	401	1/1	0.99	0.04	78,78,78,78	0
3	CA	B	401	1/1	0.99	0.03	100,100,100,100	0
3	CA	A	402	1/1	0.99	0.06	61,61,61,61	0
3	CA	B	403	1/1	0.99	0.08	74,74,74,74	0
3	CA	A	403	1/1	1.00	0.03	64,64,64,64	0
3	CA	B	404	1/1	1.00	0.02	58,58,58,58	0
3	CA	A	404	1/1	1.00	0.02	60,60,60,60	0
3	CA	A	405	1/1	1.00	0.05	52,52,52,52	0

6.5 Other polymers

There are no such residues in this entry.