



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 11:28 AM BST

PDB ID : 5DZX  
Title : Protocadherin beta 6 extracellular cadherin domains 1-4  
Authors : Goodman, K.M.; Mannepalli, S.; Bahna, F.; Honig, B.; Shapiro, L.  
Deposited on : 2015-09-26  
Resolution : 2.88 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](mailto: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.

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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

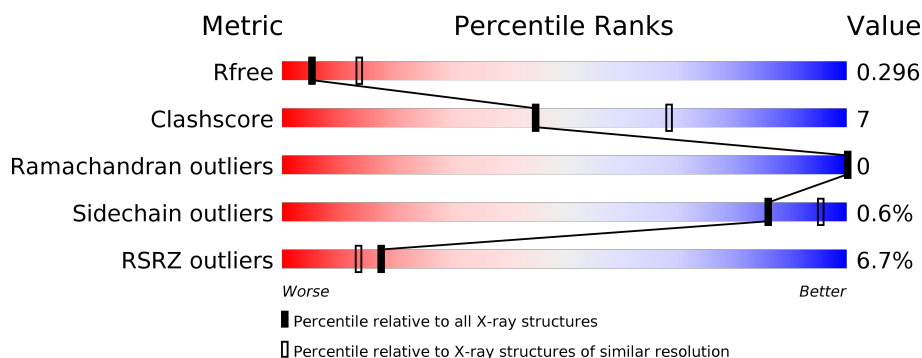
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.88 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	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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  $\geq 3$ , 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	425	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>20%</div> <div></div> </div> <div></div> </div>
1	B	425	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div></div> </div> <div></div> </div>
2	C	5	<div> <div></div> <div> <div></div> <div>80%</div> <div>20%</div> <div></div> </div> <div></div> </div>
2	D	5	<div> <div></div> <div> <div></div> <div>60%</div> <div>40%</div> <div></div> </div> <div></div> </div>
3	E	2	<div> <div></div> <div> <div></div> <div>100%</div> <div></div> </div> <div></div> </div>

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	MAN	D	4	-	-	-	X

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6413 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.

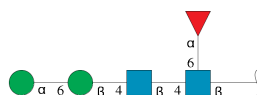
- Molecule 1 is a protein called Protocadherin beta 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	1	0
			3138	1978	523	630	7			
1	B	411	Total	C	N	O	S	0	0	0
			3069	1936	508	618	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	418	HIS	-	expression tag	UNP B2RWV0
A	419	HIS	-	expression tag	UNP B2RWV0
A	420	HIS	-	expression tag	UNP B2RWV0
A	421	HIS	-	expression tag	UNP B2RWV0
A	422	HIS	-	expression tag	UNP B2RWV0
A	423	HIS	-	expression tag	UNP B2RWV0
A	424	HIS	-	expression tag	UNP B2RWV0
A	425	HIS	-	expression tag	UNP B2RWV0
B	418	HIS	-	expression tag	UNP B2RWV0
B	419	HIS	-	expression tag	UNP B2RWV0
B	420	HIS	-	expression tag	UNP B2RWV0
B	421	HIS	-	expression tag	UNP B2RWV0
B	422	HIS	-	expression tag	UNP B2RWV0
B	423	HIS	-	expression tag	UNP B2RWV0
B	424	HIS	-	expression tag	UNP B2RWV0
B	425	HIS	-	expression tag	UNP B2RWV0

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



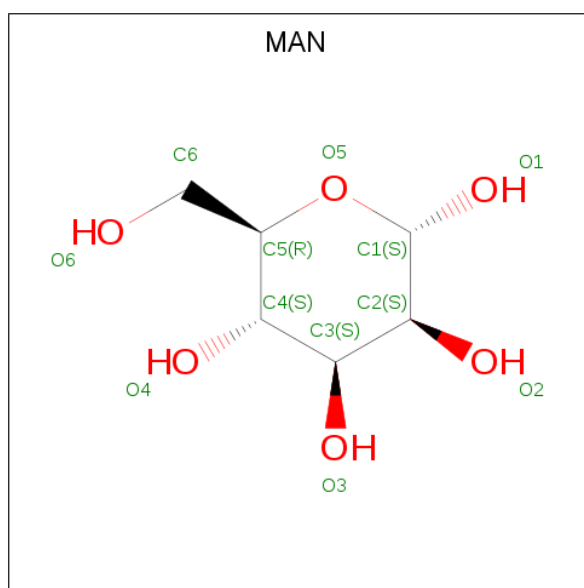
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	5	Total	C	N	O	0	0	0
			60	34	2	24			
2	D	5	Total	C	N	O	0	0	0
			60	34	2	24			

- Molecule 3 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
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is alpha-D-mannopyranose (three-letter code: MAN) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
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 CALCIUM ION (three-letter code: CA) (formula: Ca).

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

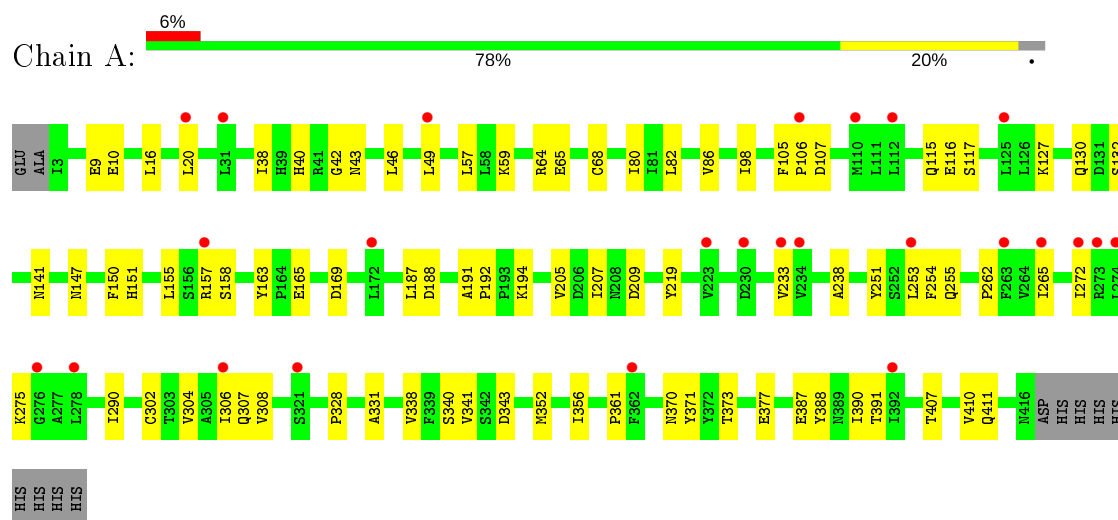
- Molecule 6 is water.

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

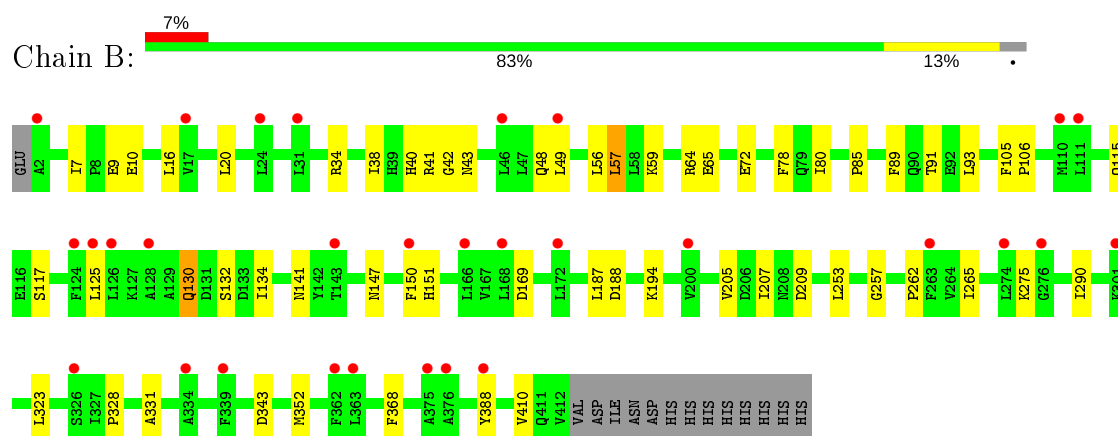
### 3 Residue-property plots [i](#)

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.

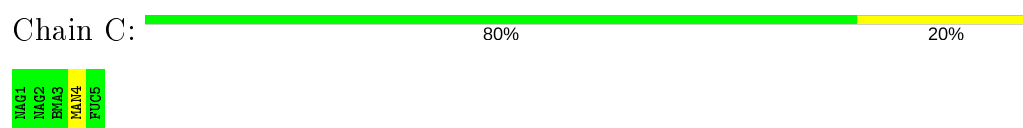
- Molecule 1: Protocadherin beta 6



- Molecule 1: Protocadherin beta 6



- Molecule 2: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2:  $\alpha$ -D-mannopyranose-(1-6)- $\beta$ -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- $\beta$ -D-glucopyranose-(1-4)-[ $\alpha$ -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- $\beta$ -D-glucopyranose

Chain D:  60% 40%

MAG1  
MAG2  
EUA3  
MAG4  
FUC5

- Molecule 3: 2-acetamido-2-deoxy- $\beta$ -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- $\beta$ -D-glucopyranose

Chain E:  100%

MAG1  
MAG2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.04Å 90.53Å 258.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.97 – 2.88 85.43 – 2.88	Depositor EDS
% Data completeness (in resolution range)	54.9 (29.97-2.88) 55.1 (85.43-2.88)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.55 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.241 , 0.289 0.247 , 0.296	Depositor DCC
$R_{free}$ test set	753 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.7	Xtriage
Anisotropy	1.198	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 43.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6413	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.63 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6981e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MAN, CA, BMA, NAG, FUC

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	0.23	0/3199	0.42	0/4376
1	B	0.22	0/3126	0.41	0/4280
All	All	0.22	0/6325	0.41	0/8656

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	3138	0	3051	51	0
1	B	3069	0	2962	35	0
2	C	60	0	52	0	0
2	D	60	0	52	1	0
3	E	28	0	25	0	0
4	A	11	0	10	0	0
4	B	22	0	20	0	0
5	A	9	0	0	0	0
5	B	9	0	0	0	0
6	A	3	0	0	0	0
6	B	4	0	0	0	0
All	All	6413	0	6172	82	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 (82) 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:9:GLU:HG2	1:B:10:GLU:HG2	1.75	0.67
1:A:9:GLU:HG2	1:A:10:GLU:HG2	1.78	0.66
1:A:20:LEU:HD12	1:A:49:LEU:HD13	1.78	0.65
1:A:46:LEU:O	1:A:59:LYS:N	2.28	0.65
1:A:147:ASN:HD21	1:A:150:PHE:HB2	1.62	0.63
1:A:107:ASP:OD2	1:A:127:LYS:NZ	2.27	0.63
1:B:147:ASN:HD21	1:B:150:PHE:HB2	1.64	0.62
1:B:9:GLU:HG3	1:B:64:ARG:H	1.62	0.62
1:A:151:HIS:ND1	1:A:169:ASP:OD2	2.33	0.61
1:A:219:TYR:HB2	1:A:304:VAL:HG22	1.83	0.61
1:A:9:GLU:HG3	1:A:64:ARG:H	1.64	0.61
1:A:157:ARG:HG2	1:A:158:SER:H	1.66	0.61
1:A:38:ILE:HG13	1:A:80:ILE:HG22	1.82	0.60
1:B:115:GLN:HA	1:B:205:VAL:HB	1.82	0.60
1:B:34:ARG:NH2	1:B:85:PRO:O	2.35	0.59
1:A:157:ARG:HB3	1:A:163:TYR:HE2	1.67	0.59
1:A:157:ARG:HB3	1:A:163:TYR:CE2	2.38	0.58
1:B:265:ILE:HD12	1:B:290:ILE:HD12	1.85	0.58
1:A:65:GLU:HG2	1:A:132:SER:HB3	1.87	0.55
1:A:265:ILE:HD13	1:A:272:ILE:HG12	1.90	0.54
1:A:40:HIS:NE2	1:A:43:ASN:OD1	2.41	0.54
1:A:391:THR:HG21	2:D:1:NAG:H5	1.90	0.54
1:B:65:GLU:HG2	1:B:132:SER:HB3	1.92	0.52
1:A:253:LEU:HD23	1:A:265:ILE:HB	1.91	0.52
1:B:20:LEU:HD12	1:B:49:LEU:HD13	1.92	0.51
1:B:188:ASP:HB3	1:B:194:LYS:H	1.76	0.51
1:A:388:TYR:HB2	1:A:410:VAL:HG13	1.91	0.51
1:A:40:HIS:HD2	1:A:42:GLY:H	1.59	0.50
1:B:141:ASN:HB3	1:B:187:LEU:HB2	1.94	0.50
1:B:48:GLN:HB2	1:B:59:LYS:HG3	1.93	0.50
1:A:141:ASN:HB3	1:A:187:LEU:HB2	1.93	0.49
1:A:341:VAL:HG23	1:A:370:ASN:HB3	1.94	0.49
1:A:238:ALA:N	1:A:251:TYR:OH	2.44	0.49
1:A:343:ASP:HB2	1:A:352:MET:SD	2.52	0.49
1:B:262:PRO:O	1:B:275:LYS:N	2.35	0.49
1:A:262:PRO:O	1:A:275:LYS:N	2.36	0.49
1:A:387:GLU:OE2	1:A:411:GLN:NE2	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:PRO:HB2	1:B:331:ALA:HB2	1.94	0.48
1:B:16:LEU:HA	1:B:57:LEU:HB3	1.95	0.48
1:B:20:LEU:HD11	1:B:56:LEU:HD22	1.95	0.48
1:A:338:VAL:HG22	1:A:373:THR:HG22	1.97	0.47
1:A:188:ASP:HB3	1:A:194:LYS:H	1.79	0.47
1:B:117:SER:HB3	1:B:207:ILE:HD13	1.97	0.46
1:A:130:GLN:HA	1:B:257:GLY:O	2.16	0.46
1:B:388:TYR:HB2	1:B:410:VAL:HG13	1.98	0.46
1:A:117:SER:HB3	1:A:207:ILE:HD13	1.97	0.46
1:A:307[B]:GLN:NE2	1:A:308:VAL:O	2.40	0.46
1:B:43:ASN:N	1:B:43:ASN:OD1	2.48	0.46
1:A:116:GLU:HB3	1:A:207:ILE:HG12	1.96	0.45
1:A:115:GLN:HA	1:A:205:VAL:HB	1.98	0.45
1:A:341:VAL:O	1:B:41:ARG:NH2	2.49	0.45
1:A:290:ILE:HB	1:A:302:CYS:SG	2.56	0.45
1:A:255:GLN:NE2	1:B:130:GLN:OE1	2.49	0.45
1:A:340:SER:HB3	1:A:371:TYR:CE2	2.52	0.45
1:A:233:VAL:HG11	1:A:306:ILE:HD13	1.99	0.45
1:B:7:ILE:HG12	1:B:93:LEU:HD11	1.98	0.44
1:A:209:ASP:N	1:A:209:ASP:OD1	2.50	0.44
1:B:253:LEU:HD23	1:B:265:ILE:HB	1.99	0.44
1:A:105:PHE:HA	1:A:106:PRO:HD3	1.90	0.44
1:B:80:ILE:HG13	1:B:89:PHE:HB2	1.99	0.44
1:A:16:LEU:HA	1:A:57:LEU:HG	1.99	0.43
1:B:38:ILE:HG13	1:B:80:ILE:HG22	1.98	0.43
1:B:78:PHE:HB2	1:B:91:THR:HG23	1.99	0.43
1:B:64:ARG:HH21	1:B:65:GLU:HG3	1.83	0.43
1:A:191:ALA:HA	1:A:192:PRO:HA	1.91	0.42
1:B:105:PHE:HA	1:B:106:PRO:HD3	1.88	0.42
1:B:40:HIS:HD2	1:B:42:GLY:H	1.66	0.42
1:A:9:GLU:HB3	1:A:98:ILE:HG13	2.02	0.42
1:A:361:PRO:O	1:A:377:GLU:HB3	2.20	0.42
1:A:64:ARG:HD2	1:A:68:CYS:HB2	2.00	0.42
1:B:9:GLU:HG2	1:B:10:GLU:N	2.35	0.42
1:A:9:GLU:HG2	1:A:10:GLU:N	2.35	0.41
1:B:343:ASP:HB2	1:B:352:MET:SD	2.59	0.41
1:A:254:PHE:CD2	1:B:125:LEU:HD12	2.56	0.41
1:B:151:HIS:ND1	1:B:169:ASP:OD2	2.50	0.41
1:B:72:GLU:HG3	1:B:134:ILE:HD13	2.02	0.41
1:B:209:ASP:OD1	1:B:209:ASP:N	2.53	0.41
1:A:82:LEU:O	1:A:86:VAL:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LEU:HD12	1:A:165:GLU:HG3	2.02	0.41
1:A:391:THR:HG22	1:A:407:THR:HB	2.03	0.41
1:A:356:ILE:HD11	1:A:390:ILE:HG23	2.03	0.40
1:A:328:PRO:HB2	1:A:331:ALA:HB2	2.04	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	413/425 (97%)	393 (95%)	20 (5%)	0	100	100
1	B	409/425 (96%)	389 (95%)	20 (5%)	0	100	100
All	All	822/850 (97%)	782 (95%)	40 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 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	346/370 (94%)	346 (100%)	0	100	100
1	B	334/370 (90%)	330 (99%)	4 (1%)	71	89
All	All	680/740 (92%)	676 (99%)	4 (1%)	86	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	57	LEU
1	B	130	GLN
1	B	323	LEU
1	B	368	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	ASN
1	B	147	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 ⓘ

12 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	1,2	14,14,15	0.48	0	17,19,21	0.58	0
2	NAG	C	2	2	14,14,15	0.47	0	17,19,21	0.46	0
2	BMA	C	3	2	11,11,12	0.62	0	15,15,17	0.94	0
2	MAN	C	4	2	11,11,12	0.92	1 (9%)	15,15,17	0.90	1 (6%)
2	FUC	C	5	2	10,10,11	0.83	0	14,14,16	0.85	0
2	NAG	D	1	1,2	14,14,15	0.43	0	17,19,21	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	D	2	2	14,14,15	0.32	0	17,19,21	0.43	0
2	BMA	D	3	2	11,11,12	0.66	0	15,15,17	0.73	0
2	MAN	D	4	2	11,11,12	0.73	0	15,15,17	1.15	2 (13%)
2	FUC	D	5	2	10,10,11	0.79	0	14,14,16	0.83	0
3	NAG	E	1	1,3	14,14,15	0.35	0	17,19,21	0.69	1 (5%)
3	NAG	E	2	3	14,14,15	0.56	0	17,19,21	0.67	1 (5%)

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	1,2	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	2/2/19/22	0/1/1/1
2	MAN	C	4	2	-	1/2/19/22	0/1/1/1
2	FUC	C	5	2	-	-	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	BMA	D	3	2	-	1/2/19/22	0/1/1/1
2	MAN	D	4	2	-	0/2/19/22	0/1/1/1
2	FUC	D	5	2	-	-	0/1/1/1
3	NAG	E	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	MAN	O5-C1	-2.38	1.39	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	MAN	C1-O5-C5	2.70	115.85	112.19
3	E	2	NAG	C1-O5-C5	2.53	115.63	112.19
2	D	4	MAN	O2-C2-C3	-2.19	105.75	110.14
2	C	4	MAN	O2-C2-C3	-2.18	105.78	110.14
3	E	1	NAG	C1-O5-C5	2.09	115.03	112.19

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	C	3	BMA	C4-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	C	1	NAG	C4-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
2	C	3	BMA	O5-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	4	MAN	O5-C5-C6-O6
2	D	3	BMA	C4-C5-C6-O6

There are no ring outliers.

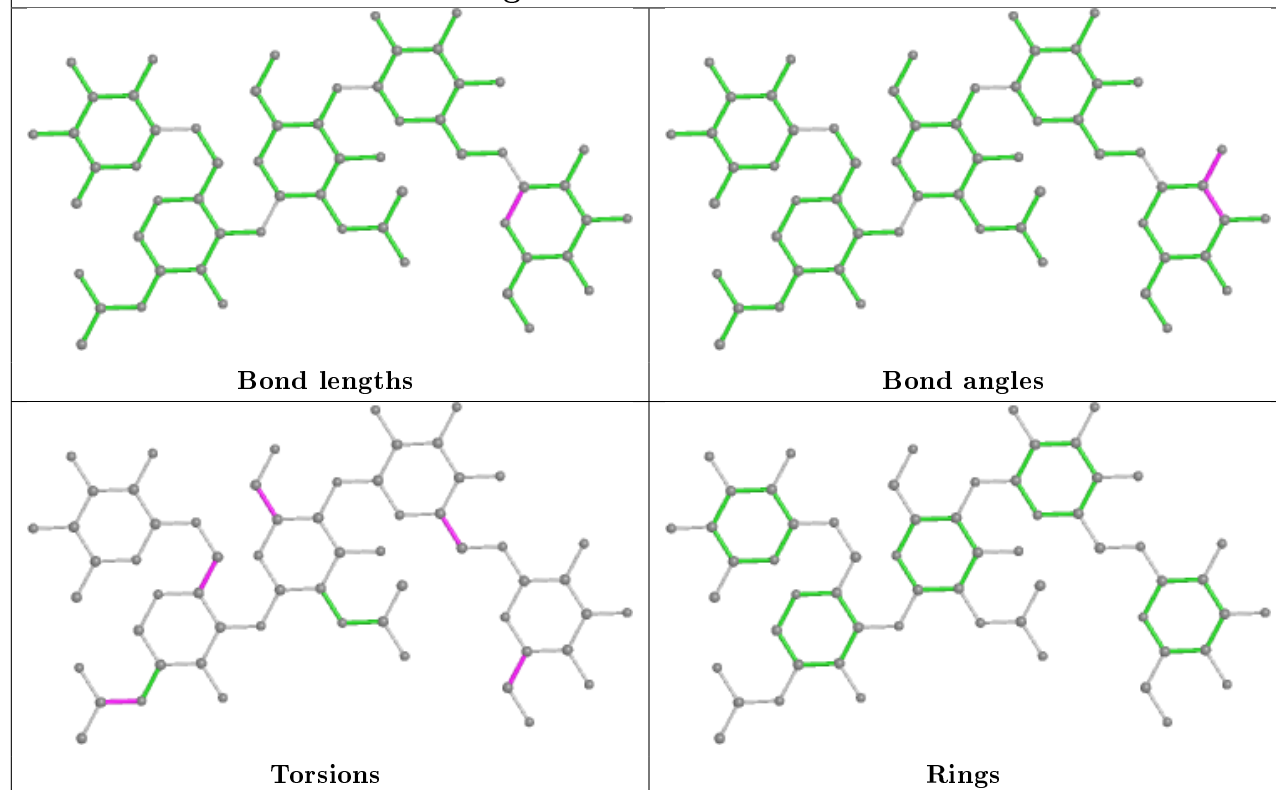
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NAG	1	0

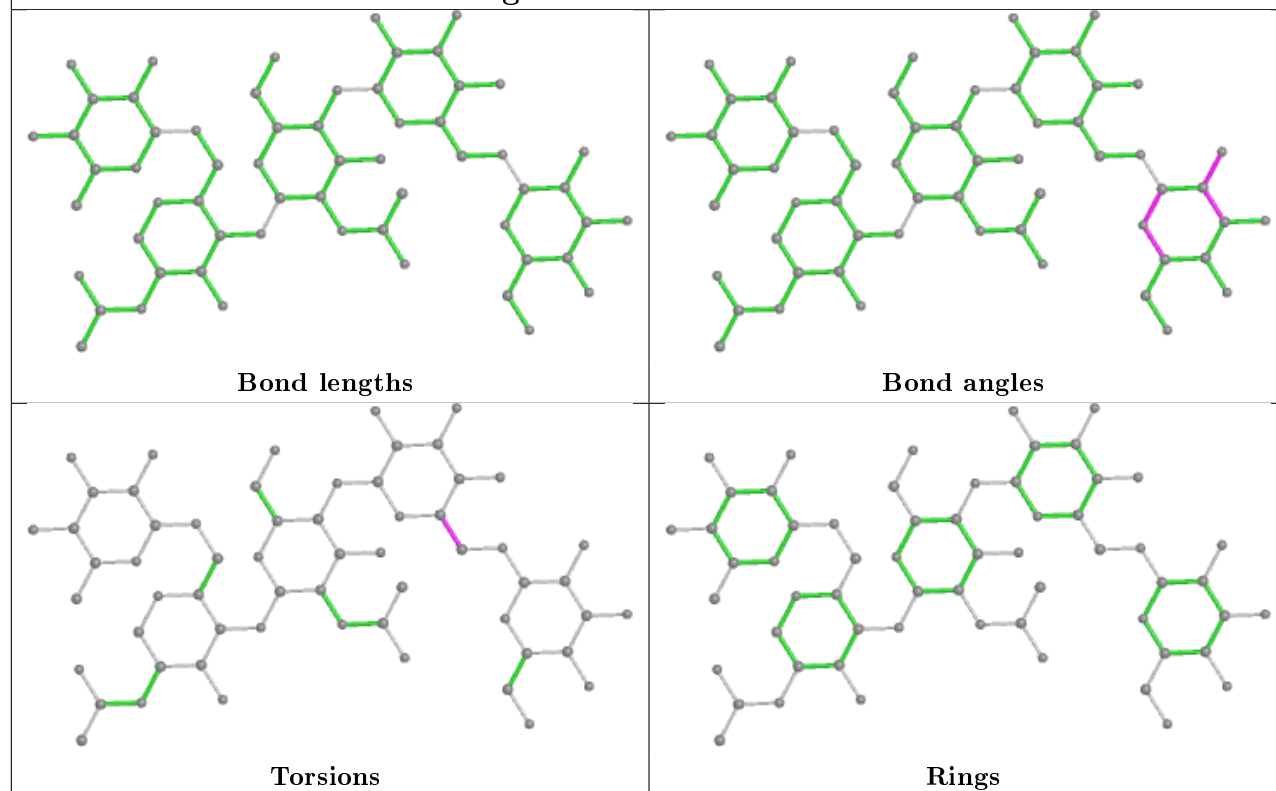
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

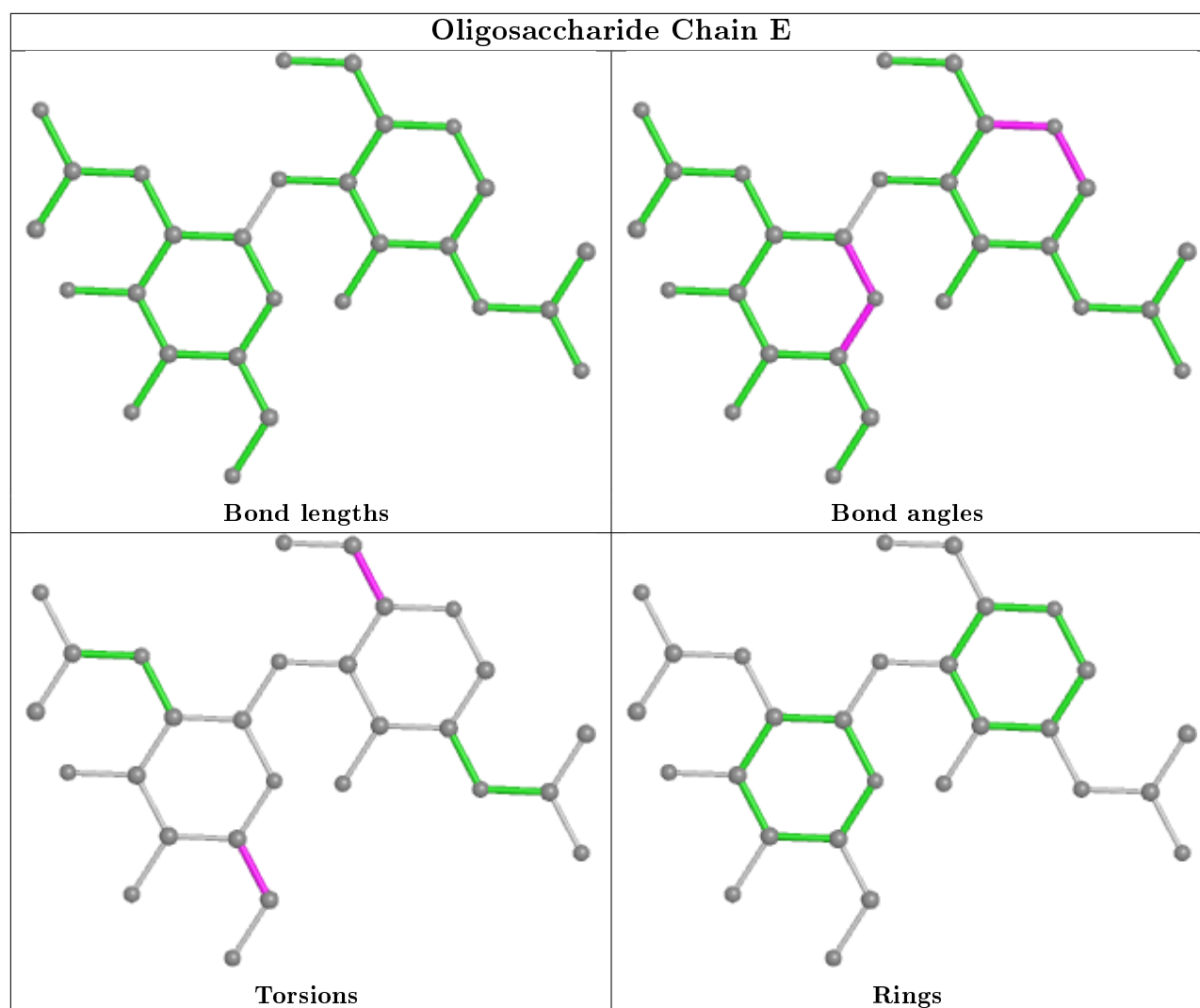


## Oligosaccharide Chain C



## Oligosaccharide Chain D





## 5.6 Ligand geometry ⓘ

Of 21 ligands modelled in this entry, 18 are monoatomic - leaving 3 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	B	504	1	11,11,12	0.72	0	15,15,17	1.08	2 (13%)
4	MAN	A	511	1	11,11,12	0.69	0	15,15,17	1.15	2 (13%)
4	MAN	B	503	1	11,11,12	0.65	0	15,15,17	1.21	1 (6%)

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	B	504	1	-	0/2/19/22	0/1/1/1
4	MAN	A	511	1	-	0/2/19/22	0/1/1/1
4	MAN	B	503	1	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	503	MAN	C1-O5-C5	3.28	116.64	112.19
4	A	511	MAN	C1-O5-C5	3.06	116.34	112.19
4	B	504	MAN	C1-O5-C5	2.74	115.90	112.19
4	A	511	MAN	O2-C2-C3	-2.17	105.79	110.14
4	B	504	MAN	O2-C2-C3	-2.12	105.89	110.14

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	414/425 (97%)	0.35	25 (6%) 21 17	38, 61, 88, 106	0
1	B	411/425 (96%)	0.43	30 (7%) 15 11	39, 63, 115, 134	0
All	All	825/850 (97%)	0.39	55 (6%) 17 13	38, 63, 99, 134	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	126	LEU	5.3
1	B	375	ALA	4.9
1	B	363	LEU	4.7
1	B	111	LEU	3.6
1	A	272	ILE	3.5
1	B	326	SER	3.4
1	B	46	LEU	3.3
1	B	110	MET	3.1
1	B	17	VAL	3.1
1	A	274	LEU	3.0
1	A	223	VAL	2.9
1	A	112	LEU	2.9
1	A	110	MET	2.8
1	A	392	ILE	2.8
1	B	339	PHE	2.8
1	B	125	LEU	2.8
1	B	31	LEU	2.7
1	B	172	LEU	2.7
1	A	49	LEU	2.6
1	B	166	LEU	2.6
1	A	172	LEU	2.5
1	B	124	PHE	2.5
1	B	49	LEU	2.4
1	A	233	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	265	ILE	2.4
1	A	31	LEU	2.4
1	B	388	TYR	2.4
1	B	334	ALA	2.4
1	B	276	GLY	2.3
1	B	150	PHE	2.3
1	A	362	PHE	2.3
1	A	106	PRO	2.3
1	A	230	ASP	2.3
1	A	321	SER	2.3
1	B	128	ALA	2.3
1	B	143	THR	2.3
1	A	20	LEU	2.3
1	B	263	PHE	2.3
1	A	306	ILE	2.2
1	B	274	LEU	2.2
1	B	376	ALA	2.2
1	A	125	LEU	2.2
1	A	278	LEU	2.2
1	A	263	PHE	2.1
1	A	234	VAL	2.1
1	A	276	GLY	2.1
1	B	2	ALA	2.1
1	B	24	LEU	2.1
1	B	168	LEU	2.1
1	B	362	PHE	2.1
1	B	301	LYS	2.1
1	B	200	VAL	2.1
1	A	273	ARG	2.0
1	A	157	ARG	2.0
1	A	253	LEU	2.0

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

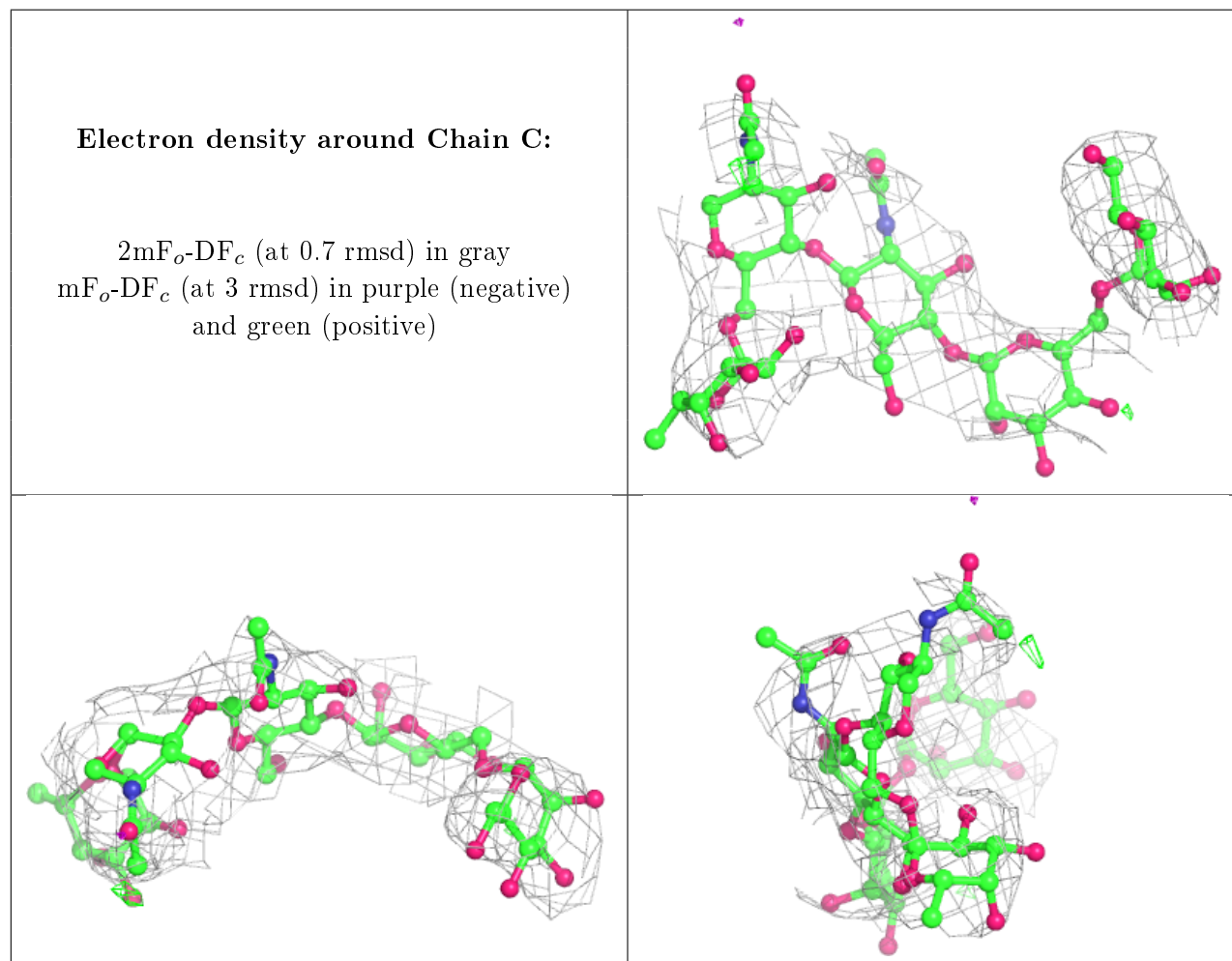
There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 95<sup>th</sup> 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.

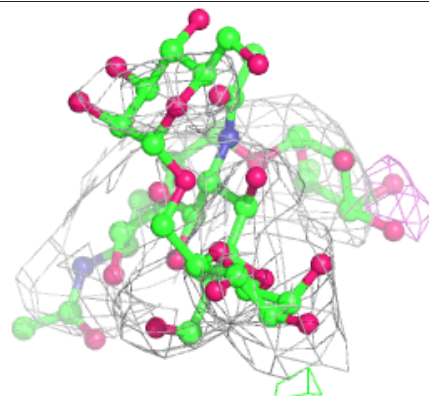
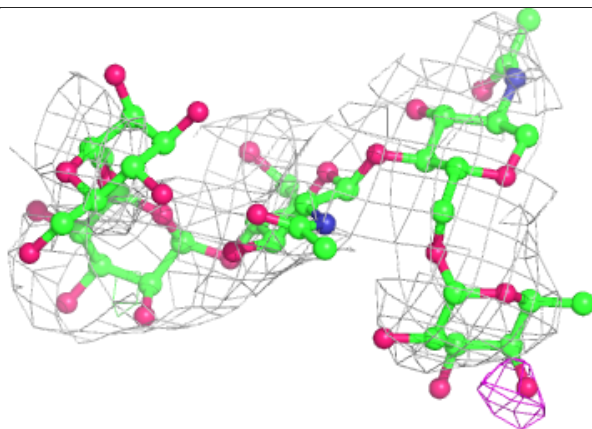
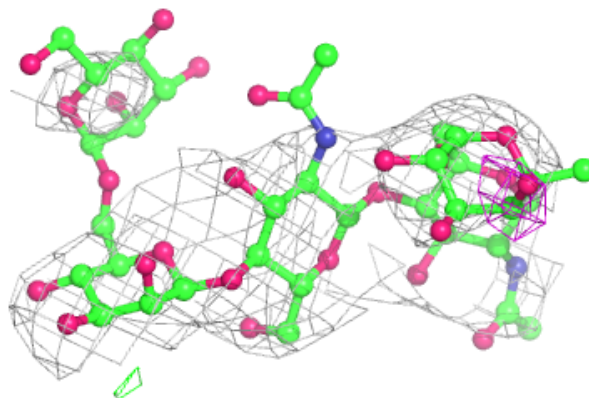
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	BMA	C	3	11/12	0.55	0.23	135,147,152,152	0
2	MAN	D	4	11/12	0.61	0.42	90,129,145,145	0
2	BMA	D	3	11/12	0.62	0.24	101,114,137,138	0
2	MAN	C	4	11/12	0.69	0.21	116,134,144,146	0
2	FUC	C	5	10/11	0.77	0.27	109,119,122,124	0
3	NAG	E	1	14/15	0.77	0.16	89,106,112,114	0
2	NAG	D	2	14/15	0.78	0.27	103,116,125,131	0
2	FUC	D	5	10/11	0.85	0.37	74,95,100,105	0
3	NAG	E	2	14/15	0.87	0.17	91,108,116,116	0
2	NAG	C	2	14/15	0.88	0.13	120,137,148,148	0
2	NAG	C	1	14/15	0.92	0.14	81,108,121,137	0
2	NAG	D	1	14/15	0.95	0.16	88,96,112,113	0

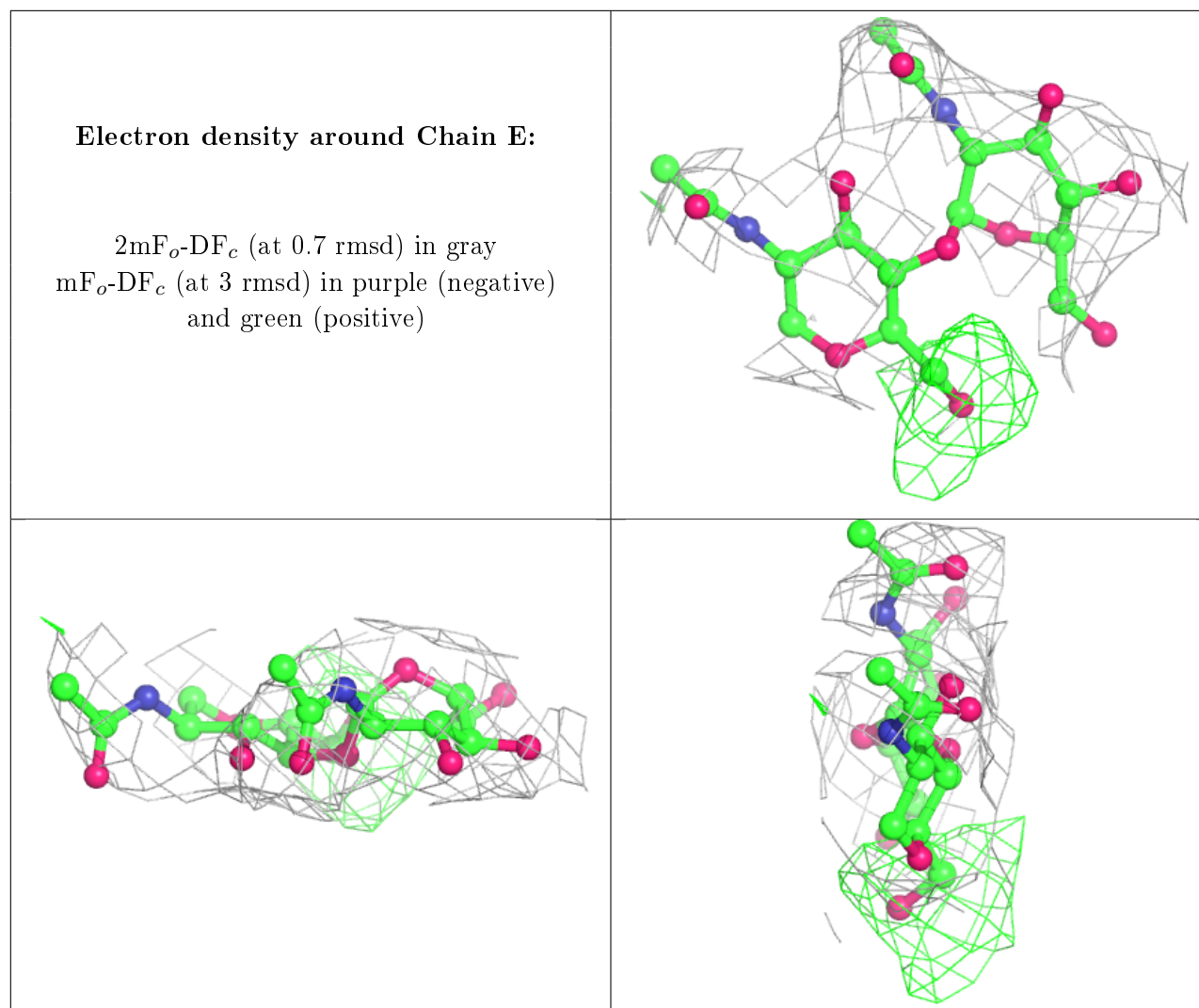
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around Chain 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	MAN	B	504	11/12	0.81	0.24	86,98,111,112	0
5	CA	B	511	1/1	0.82	0.28	63,63,63,63	0
4	MAN	B	503	11/12	0.85	0.33	79,98,109,114	0
4	MAN	A	511	11/12	0.90	0.34	70,89,114,119	0
5	CA	A	518	1/1	0.91	0.27	62,62,62,62	0
5	CA	B	507	1/1	0.94	0.19	57,57,57,57	0
5	CA	A	514	1/1	0.96	0.29	36,36,36,36	0
5	CA	A	517	1/1	0.98	0.20	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	CA	B	506	1/1	0.98	0.25	37,37,37,37	0
5	CA	B	512	1/1	0.98	0.25	22,22,22,22	0
5	CA	B	508	1/1	0.98	0.19	56,56,56,56	0
5	CA	A	516	1/1	0.98	0.23	39,39,39,39	0
5	CA	B	509	1/1	0.98	0.21	38,38,38,38	0
5	CA	A	513	1/1	0.99	0.27	39,39,39,39	0
5	CA	B	513	1/1	0.99	0.25	27,27,27,27	0
5	CA	A	519	1/1	0.99	0.20	27,27,27,27	0
5	CA	A	512	1/1	0.99	0.24	63,63,63,63	0
5	CA	B	505	1/1	0.99	0.22	53,53,53,53	0
5	CA	B	510	1/1	0.99	0.18	55,55,55,55	0
5	CA	A	515	1/1	0.99	0.18	47,47,47,47	0
5	CA	A	520	1/1	0.99	0.27	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.