



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 07:20 PM BST

PDB ID : 5DZY  
Title : Protocadherin beta 8 extracellular cadherin domains 1-4  
Authors : Goodman, K.M.; Bahna, F.; Mannepalli, S.; Honig, B.; Shapiro, L.  
Deposited on : 2015-09-26  
Resolution : 2.90 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

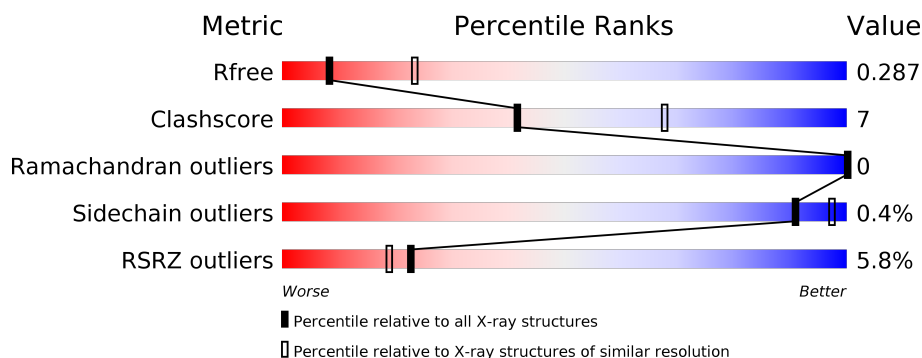
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	426	<div> <div>3%</div> <div>81%</div> <div>16%</div> <div>.</div> </div>
1	B	426	<div> <div>5%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
1	C	426	<div> <div>10%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
1	D	426	<div> <div>8%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
1	E	426	<div> <div>%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
1	F	426	<div> <div>5%</div> <div>83%</div> <div>14%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	H	2	 100%
2	K	2	 100%
3	I	9	 67%33%
4	J	3	 67%33%
5	L	5	 80%20%

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 19132 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 Pcdhb8 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3100	1948	518	627	7			
1	B	415	Total	C	N	O	S	0	0	0
			3091	1946	505	632	8			
1	C	415	Total	C	N	O	S	0	0	0
			3079	1929	516	626	8			
1	D	412	Total	C	N	O	S	0	0	0
			3055	1928	502	617	8			
1	E	413	Total	C	N	O	S	0	0	0
			3101	1953	516	624	8			
1	F	415	Total	C	N	O	S	0	0	0
			3125	1967	516	634	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	419	HIS	-	expression tag	UNP A8E4K6
A	420	HIS	-	expression tag	UNP A8E4K6
A	421	HIS	-	expression tag	UNP A8E4K6
A	422	HIS	-	expression tag	UNP A8E4K6
A	423	HIS	-	expression tag	UNP A8E4K6
A	424	HIS	-	expression tag	UNP A8E4K6
A	425	HIS	-	expression tag	UNP A8E4K6
A	426	HIS	-	expression tag	UNP A8E4K6
B	419	HIS	-	expression tag	UNP A8E4K6
B	420	HIS	-	expression tag	UNP A8E4K6
B	421	HIS	-	expression tag	UNP A8E4K6
B	422	HIS	-	expression tag	UNP A8E4K6
B	423	HIS	-	expression tag	UNP A8E4K6
B	424	HIS	-	expression tag	UNP A8E4K6
B	425	HIS	-	expression tag	UNP A8E4K6
B	426	HIS	-	expression tag	UNP A8E4K6
C	419	HIS	-	expression tag	UNP A8E4K6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	420	HIS	-	expression tag	UNP A8E4K6
C	421	HIS	-	expression tag	UNP A8E4K6
C	422	HIS	-	expression tag	UNP A8E4K6
C	423	HIS	-	expression tag	UNP A8E4K6
C	424	HIS	-	expression tag	UNP A8E4K6
C	425	HIS	-	expression tag	UNP A8E4K6
C	426	HIS	-	expression tag	UNP A8E4K6
D	419	HIS	-	expression tag	UNP A8E4K6
D	420	HIS	-	expression tag	UNP A8E4K6
D	421	HIS	-	expression tag	UNP A8E4K6
D	422	HIS	-	expression tag	UNP A8E4K6
D	423	HIS	-	expression tag	UNP A8E4K6
D	424	HIS	-	expression tag	UNP A8E4K6
D	425	HIS	-	expression tag	UNP A8E4K6
D	426	HIS	-	expression tag	UNP A8E4K6
E	419	HIS	-	expression tag	UNP A8E4K6
E	420	HIS	-	expression tag	UNP A8E4K6
E	421	HIS	-	expression tag	UNP A8E4K6
E	422	HIS	-	expression tag	UNP A8E4K6
E	423	HIS	-	expression tag	UNP A8E4K6
E	424	HIS	-	expression tag	UNP A8E4K6
E	425	HIS	-	expression tag	UNP A8E4K6
E	426	HIS	-	expression tag	UNP A8E4K6
F	419	HIS	-	expression tag	UNP A8E4K6
F	420	HIS	-	expression tag	UNP A8E4K6
F	421	HIS	-	expression tag	UNP A8E4K6
F	422	HIS	-	expression tag	UNP A8E4K6
F	423	HIS	-	expression tag	UNP A8E4K6
F	424	HIS	-	expression tag	UNP A8E4K6
F	425	HIS	-	expression tag	UNP A8E4K6
F	426	HIS	-	expression tag	UNP A8E4K6

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



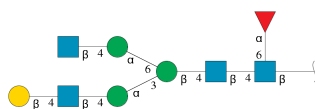
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	2	Total	C	N	O	0	0	0
			24	14	1	9			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	2	Total	C	N	O	0	0	0
			24	14	1	9			
2	K	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 3 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-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.



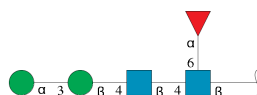
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	I	9	Total	C	N	O	0	0	0
			110	62	4	44			

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



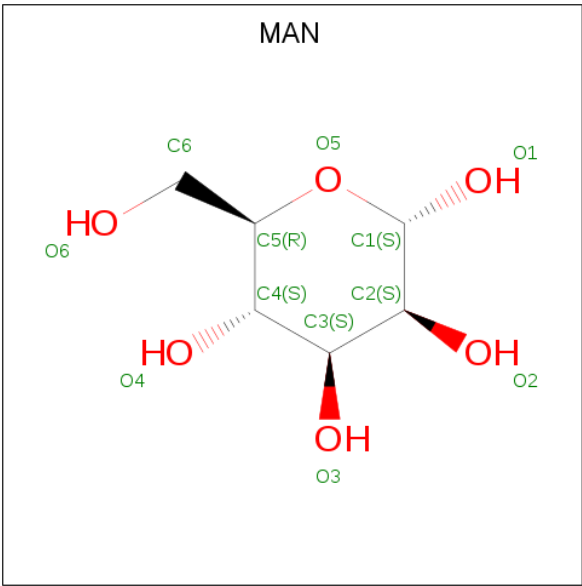
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-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
5	L	5	Total	C	N	O	0	0	0
			60	34	2	24			

- Molecule 6 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
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	A	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	B	1	Total	C	O	0	0
			11	6	5		
6	C	1	Total	C	O	0	0
			11	6	5		
6	C	1	Total	C	O	0	0
			11	6	5		
6	C	1	Total	C	O	0	0
			11	6	5		
6	D	1	Total	C	O	0	0
			11	6	5		
6	D	1	Total	C	O	0	0
			11	6	5		
6	D	1	Total	C	O	0	0
			11	6	5		
6	E	1	Total	C	O	0	0
			11	6	5		

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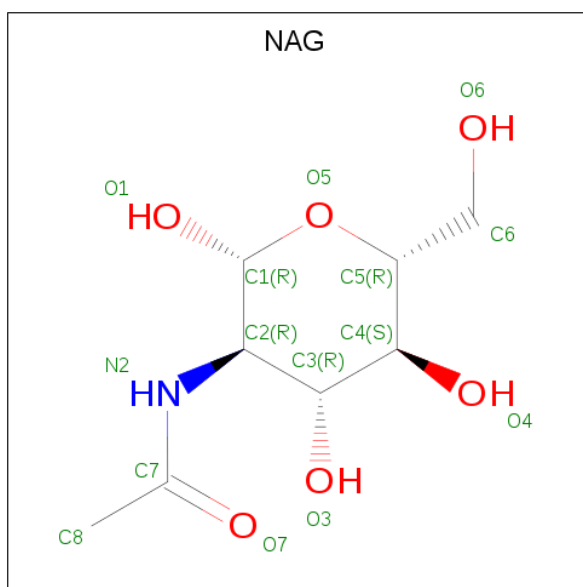
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	C	O	0	0
			11	6	5		
6	E	1	Total	C	O	0	0
			11	6	5		
6	F	1	Total	C	O	0	0
			11	6	5		
6	F	1	Total	C	O	0	0
			11	6	5		
6	F	1	Total	C	O	0	0
			11	6	5		

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	9	Total	Ca	0	0
			9	9		
7	E	9	Total	Ca	0	0
			9	9		
7	B	9	Total	Ca	0	0
			9	9		
7	C	9	Total	Ca	0	0
			9	9		
7	A	9	Total	Ca	0	0
			9	9		
7	F	9	Total	Ca	0	0
			9	9		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	1	Total	C	N	O	0	0
			14	8	1	5		
8	F	1	Total	C	N	O	0	0
			14	8	1	5		
8	F	1	Total	C	N	O	0	0
			14	8	1	5		

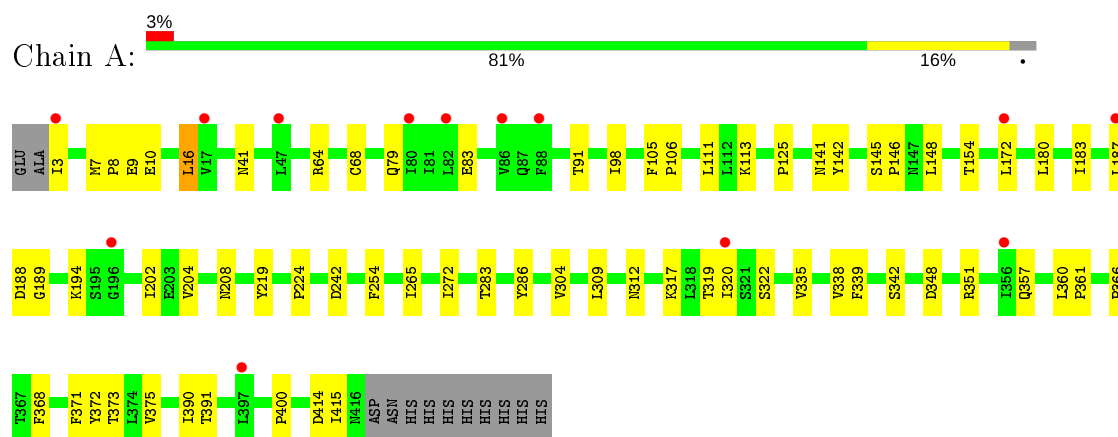
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	O	0	0
			1	1		
9	C	1	Total	O	0	0
			1	1		
9	D	1	Total	O	0	0
			1	1		
9	E	2	Total	O	0	0
			2	2		
9	F	1	Total	O	0	0
			1	1		

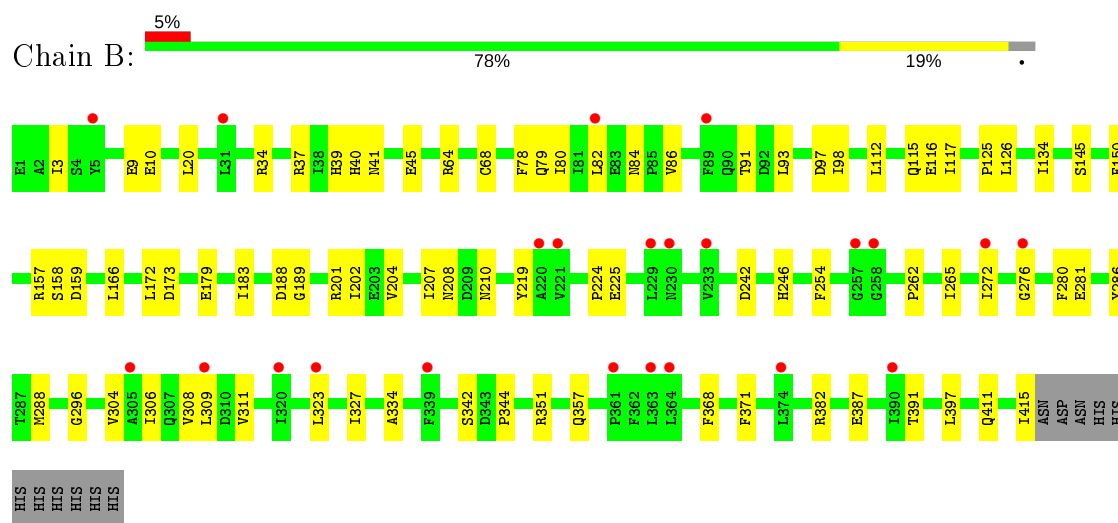
### 3 Residue-property plots [i](#)

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.

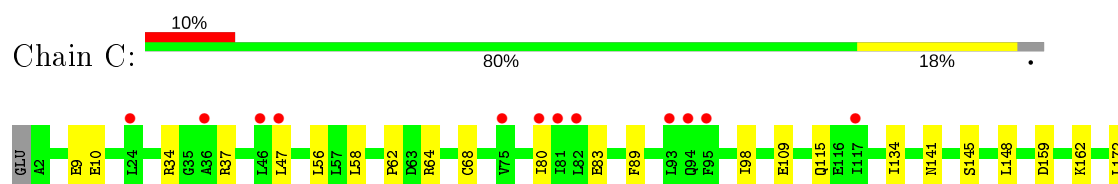
#### • Molecule 1: Pcdhb8 protein

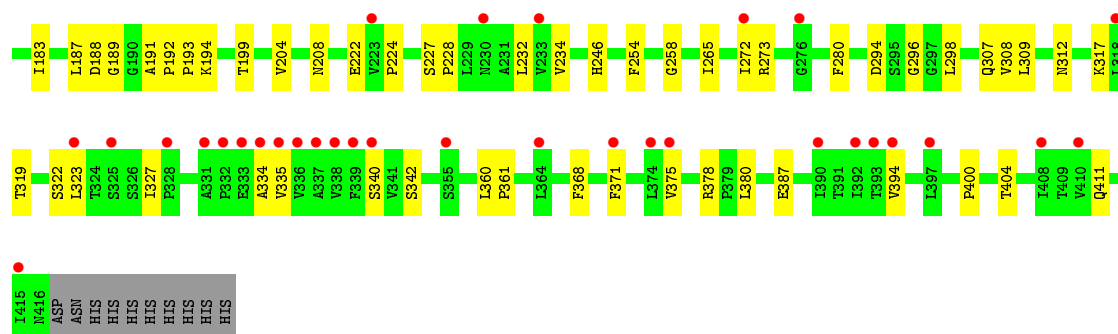


#### • Molecule 1: Pcdhb8 protein

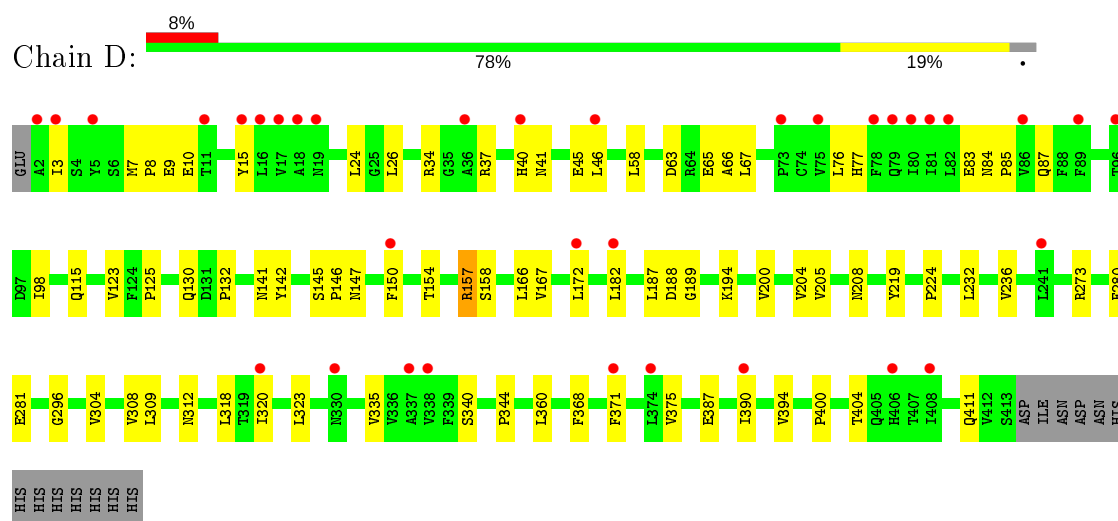


#### • Molecule 1: Pcdhb8 protein

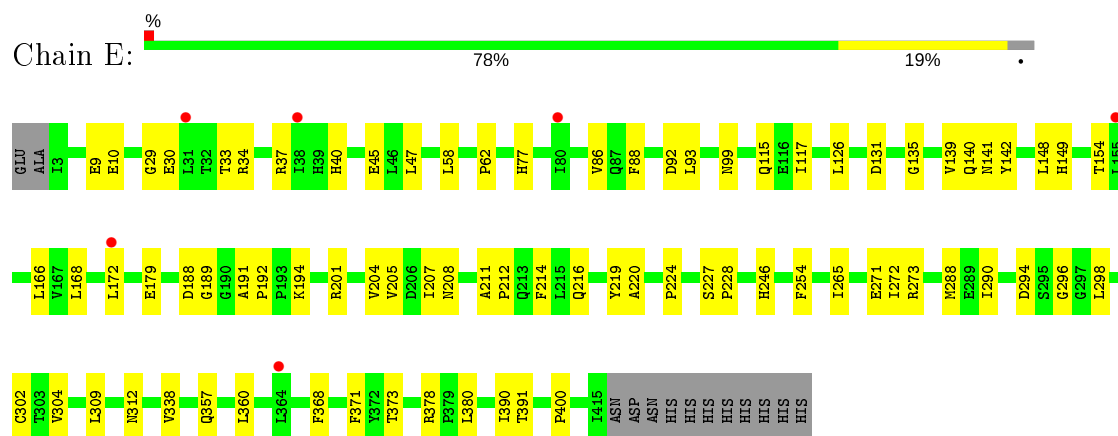




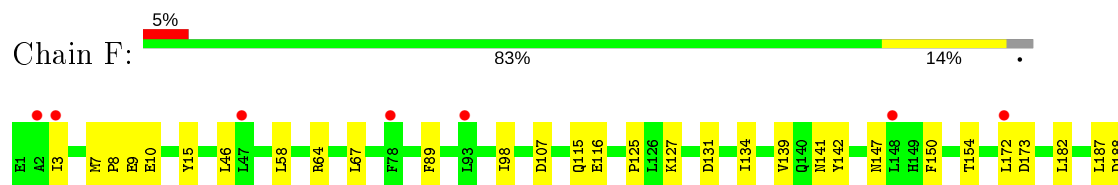
- Molecule 1: Pcdhb8 protein

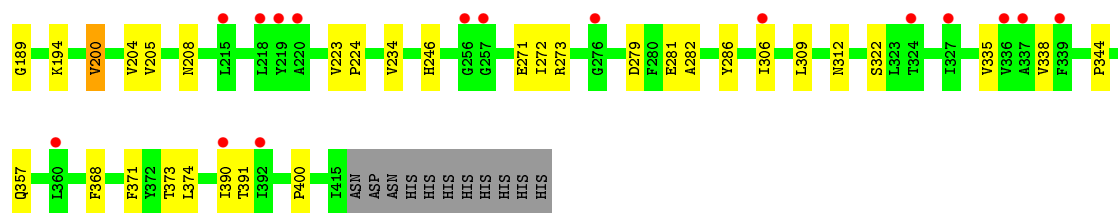


- Molecule 1: Pcdhb8 protein



- Molecule 1: Pcdhb8 protein





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

Chain G: 100%



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

Chain H: 100%



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

Chain K: 100%



- Molecule 3: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-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 I: 67% 33%



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

Chain J: 67% 33%



- Molecule 5: alpha-D-mannopyranose-(1-3)-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 L: 80% 20%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.14Å 137.07Å 326.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.99 – 2.90 126.37 – 2.72	Depositor EDS
% Data completeness (in resolution range)	62.5 (29.99-2.90) 52.3 (126.37-2.72)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.62 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.229 , 0.283 0.236 , 0.287	Depositor DCC
$R_{free}$ test set	2759 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	56.5	Xtriage
Anisotropy	0.978	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 56.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	19132	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, CA, GAL, FUC, 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	0.20	0/3158	0.38	0/4322
1	B	0.21	0/3151	0.38	0/4320
1	C	0.21	0/3137	0.38	0/4298
1	D	0.20	0/3114	0.38	0/4268
1	E	0.20	0/3161	0.38	0/4329
1	F	0.20	0/3185	0.38	0/4362
All	All	0.20	0/18906	0.38	0/25899

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	3100	0	2983	43	0
1	B	3091	0	2941	51	0
1	C	3079	0	2915	49	0
1	D	3055	0	2919	53	0
1	E	3101	0	2989	51	0
1	F	3125	0	3011	41	0
2	G	24	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	24	0	22	0	0
2	K	24	0	22	0	0
3	I	110	0	94	0	0
4	J	39	0	34	0	0
5	L	60	0	52	0	0
6	A	33	0	30	2	0
6	B	33	0	30	0	0
6	C	33	0	30	0	0
6	D	33	0	30	1	0
6	E	33	0	30	0	0
6	F	33	0	30	0	0
7	A	9	0	0	0	0
7	B	9	0	0	0	0
7	C	9	0	0	0	0
7	D	9	0	0	0	0
7	E	9	0	0	0	0
7	F	9	0	0	0	0
8	B	14	0	13	0	0
8	F	28	0	26	0	0
9	A	1	0	0	0	0
9	C	1	0	0	0	0
9	D	1	0	0	0	0
9	E	2	0	0	0	0
9	F	1	0	0	0	0
All	All	19132	0	18223	261	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 (261) 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:E:99:ASN:O	1:E:194:LYS:NZ	2.15	0.77
1:E:288:MET:HG3	1:E:304:VAL:HG13	1.69	0.75
1:B:64:ARG:HH12	1:B:134:ILE:HG12	1.49	0.75
1:C:307:GLN:HE22	1:E:216:GLN:HA	1.56	0.70
1:C:64:ARG:HH22	1:C:134:ILE:HG12	1.57	0.69
1:D:40:HIS:HE2	1:D:45:GLU:HG2	1.56	0.69
1:D:142:TYR:HB2	1:D:154:THR:HG22	1.76	0.67
1:C:368:PHE:HB2	1:C:371:PHE:HB3	1.75	0.67
1:A:254:PHE:HE2	1:B:125:PRO:HB2	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:SER:O	1:B:37:ARG:NH2	2.27	0.66
1:A:219:TYR:HB2	1:A:304:VAL:HG22	1.79	0.64
1:D:368:PHE:HB2	1:D:371:PHE:HB3	1.79	0.64
1:B:3:ILE:HG13	1:B:91:THR:HG23	1.78	0.64
1:F:141:ASN:HB3	1:F:187:LEU:HB2	1.78	0.64
1:A:348:ASP:OD1	1:A:351:ARG:NH1	2.31	0.64
1:E:37:ARG:NH2	1:F:322:SER:O	2.31	0.64
1:E:357:GLN:HE22	1:E:391:THR:H	1.46	0.63
1:F:368:PHE:HB2	1:F:371:PHE:HB3	1.80	0.63
1:A:142:TYR:HB2	1:A:154:THR:HG22	1.81	0.63
1:A:125:PRO:HB2	1:B:254:PHE:HE2	1.64	0.62
1:E:142:TYR:HB2	1:E:154:THR:HG22	1.81	0.62
1:B:368:PHE:HB2	1:B:371:PHE:HB3	1.80	0.62
1:F:357:GLN:NE2	1:F:391:THR:OG1	2.32	0.62
1:F:107:ASP:OD2	1:F:127:LYS:NZ	2.27	0.62
1:B:34:ARG:HG2	1:B:84:ASN:HB3	1.82	0.62
1:A:357:GLN:NE2	1:A:391:THR:OG1	2.32	0.61
1:E:254:PHE:HE2	1:F:125:PRO:HB2	1.66	0.61
1:B:20:LEU:HD21	1:B:80:ILE:HD13	1.83	0.61
1:F:182:LEU:HB2	1:F:200:VAL:HG13	1.82	0.61
1:F:271:GLU:OE1	1:F:273:ARG:NH2	2.33	0.60
1:E:115:GLN:HA	1:E:205:VAL:HB	1.83	0.60
1:B:219:TYR:HB2	1:B:304:VAL:HG22	1.84	0.59
1:F:147:ASN:HD21	1:F:150:PHE:HB2	1.68	0.59
1:A:188:ASP:HB3	1:A:194:LYS:H	1.68	0.58
1:D:115:GLN:HA	1:D:205:VAL:HB	1.86	0.58
1:F:172:LEU:HD13	1:F:204:VAL:HG22	1.84	0.58
1:E:368:PHE:HB2	1:E:371:PHE:HB3	1.85	0.58
1:E:224:PRO:HA	1:E:309:LEU:HB2	1.85	0.58
1:F:7:MET:HE1	1:F:58:LEU:HD11	1.86	0.58
1:A:224:PRO:HA	1:A:309:LEU:HB2	1.86	0.57
1:C:148:LEU:HD11	1:F:89:PHE:HB3	1.87	0.57
1:A:360:LEU:HD13	1:A:390:ILE:HG12	1.87	0.57
1:E:47:LEU:HD11	1:E:93:LEU:HD11	1.87	0.56
1:A:312:ASN:ND2	1:A:400:PRO:O	2.37	0.56
6:A:506:MAN:HO2	6:A:507:MAN:HO4	1.53	0.56
1:D:182:LEU:HB2	1:D:200:VAL:HG13	1.87	0.56
1:B:188:ASP:OD1	1:B:189:GLY:N	2.38	0.56
1:C:387:GLU:HG3	1:C:411:GLN:HG2	1.88	0.56
1:D:9:GLU:HB3	1:D:98:ILE:HG22	1.86	0.56
1:F:64:ARG:HH12	1:F:134:ILE:HG12	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:9:GLU:HB3	1:F:98:ILE:HG22	1.87	0.56
1:C:312:ASN:ND2	1:C:400:PRO:O	2.35	0.56
1:D:147:ASN:HD21	1:D:150:PHE:HB2	1.70	0.55
1:C:58:LEU:HD21	1:C:62:PRO:HD3	1.87	0.55
1:C:232:LEU:HA	1:C:273:ARG:HG2	1.88	0.55
1:C:141:ASN:HB3	1:C:187:LEU:HB2	1.89	0.55
1:A:16:LEU:HD23	1:A:16:LEU:H	1.70	0.55
1:A:188:ASP:OD1	1:A:189:GLY:N	2.40	0.54
1:C:378:ARG:HH22	1:C:380:LEU:HD13	1.72	0.54
1:F:115:GLN:HA	1:F:205:VAL:HB	1.90	0.54
1:C:9:GLU:HB3	1:C:98:ILE:HG22	1.88	0.54
1:C:323:LEU:HA	1:D:37:ARG:HH22	1.73	0.54
1:A:368:PHE:HB2	1:A:371:PHE:HB3	1.89	0.54
1:E:58:LEU:HD21	1:E:62:PRO:HD3	1.88	0.54
1:D:232:LEU:HA	1:D:273:ARG:HG2	1.89	0.54
1:C:319:THR:OG1	1:D:41:ASN:OD1	2.26	0.53
1:D:360:LEU:HD13	1:D:390:ILE:HG12	1.91	0.53
1:D:24:LEU:HD23	1:E:148:LEU:HD21	1.91	0.53
1:A:283:THR:HG1	1:A:286:TYR:HH	1.57	0.53
1:F:188:ASP:OD1	1:F:189:GLY:N	2.42	0.53
1:B:39:HIS:HB3	1:B:79:GLN:HG2	1.90	0.53
1:D:188:ASP:OD1	1:D:189:GLY:N	2.42	0.53
1:A:265:ILE:HD13	1:A:272:ILE:HG12	1.91	0.52
1:B:281:GLU:HG2	1:B:344:PRO:HB2	1.90	0.52
1:E:312:ASN:ND2	1:E:400:PRO:O	2.42	0.52
1:C:394:VAL:HG13	1:C:404:THR:HG23	1.92	0.52
1:C:148:LEU:HD13	1:F:3:ILE:HG21	1.91	0.52
1:F:234:VAL:HG12	1:F:272:ILE:HB	1.92	0.52
1:B:382:ARG:HD3	1:B:415:ILE:HA	1.92	0.52
1:D:24:LEU:HB3	1:D:26:LEU:HD13	1.91	0.52
1:D:34:ARG:NH1	1:D:87:GLN:OE1	2.35	0.51
1:E:188:ASP:OD1	1:E:189:GLY:N	2.44	0.51
1:E:188:ASP:HB3	1:E:194:LYS:H	1.75	0.51
1:D:172:LEU:HD13	1:D:204:VAL:HG22	1.93	0.51
1:D:394:VAL:HG13	1:D:404:THR:HG23	1.91	0.51
1:D:15:TYR:HB3	1:D:58:LEU:HD12	1.91	0.51
1:C:47:LEU:HD13	1:C:56:LEU:HD21	1.92	0.50
1:B:172:LEU:HD13	1:B:204:VAL:HG22	1.92	0.50
1:A:3:ILE:HG13	1:A:91:THR:HG23	1.93	0.50
1:C:327:ILE:HD11	1:C:334:ALA:HB1	1.93	0.50
1:F:286:TYR:HB2	1:F:306:ILE:HG23	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:335:VAL:HG22	1:C:375:VAL:HG22	1.94	0.50
1:B:145:SER:HB3	1:B:183:ILE:HB	1.94	0.50
1:B:64:ARG:NH1	1:B:97:ASP:HB2	2.27	0.50
1:C:294:ASP:OD2	1:C:298:LEU:HB2	2.12	0.50
1:E:117:ILE:HG12	1:E:207:ILE:HG13	1.94	0.50
1:B:224:PRO:HA	1:B:309:LEU:HB2	1.93	0.49
1:D:281:GLU:HG2	1:D:344:PRO:HB2	1.92	0.49
1:A:111:LEU:HD21	1:A:113:LYS:HE3	1.93	0.49
1:C:188:ASP:OD1	1:C:189:GLY:N	2.45	0.49
1:C:222:GLU:HG3	1:E:220:ALA:HB3	1.94	0.49
1:B:351:ARG:HD2	1:B:397:LEU:HD12	1.93	0.49
1:B:78:PHE:HE2	1:B:93:LEU:HB2	1.77	0.49
1:E:86:VAL:HG11	1:F:335:VAL:HB	1.94	0.49
1:B:225:GLU:HB3	1:B:311:VAL:HG22	1.93	0.49
1:C:109:GLU:HA	1:C:199:THR:HB	1.95	0.49
1:F:224:PRO:HA	1:F:309:LEU:HB2	1.95	0.49
1:D:312:ASN:ND2	1:D:400:PRO:O	2.46	0.49
1:A:317:LYS:HB3	1:A:342:SER:HB3	1.95	0.48
1:A:319:THR:HG21	1:B:41:ASN:O	2.13	0.48
1:E:172:LEU:HD13	1:E:204:VAL:HG22	1.94	0.48
1:F:281:GLU:HG2	1:F:344:PRO:HB2	1.95	0.48
1:E:77:HIS:ND1	1:E:92:ASP:OD1	2.46	0.48
1:A:338:VAL:HG23	1:A:373:THR:HG22	1.96	0.48
1:B:327:ILE:HD11	1:B:334:ALA:HB1	1.96	0.48
1:B:150:PHE:HB3	1:B:166:LEU:HD11	1.96	0.48
1:F:142:TYR:HD2	1:F:154:THR:HG23	1.78	0.48
1:D:40:HIS:NE2	1:D:45:GLU:HG2	2.28	0.48
1:C:188:ASP:HB3	1:C:194:LYS:H	1.78	0.47
1:C:340:SER:HB3	1:C:371:PHE:HD1	1.79	0.47
1:F:224:PRO:HG3	1:F:309:LEU:HD12	1.96	0.47
1:A:41:ASN:HD22	1:B:342:SER:HB2	1.80	0.47
1:B:117:ILE:HG12	1:B:207:ILE:HG13	1.96	0.47
1:D:34:ARG:O	1:D:83:GLU:N	2.39	0.47
1:E:360:LEU:HD13	1:E:390:ILE:HG12	1.96	0.47
1:A:83:GLU:HG2	1:B:323:LEU:HD22	1.97	0.47
1:B:158:SER:OG	1:B:159:ASP:N	2.46	0.47
1:D:150:PHE:HB3	1:D:166:LEU:HD11	1.96	0.47
1:C:234:VAL:HG12	1:C:272:ILE:HB	1.96	0.46
1:D:46:LEU:HD21	1:D:67:LEU:HD21	1.96	0.46
1:A:172:LEU:HD13	1:A:204:VAL:HG22	1.97	0.46
1:C:37:ARG:NH2	1:D:323:LEU:HA	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:223:VAL:HB	1:F:306:ILE:HD11	1.96	0.46
1:A:335:VAL:HG22	1:A:375:VAL:HG22	1.97	0.46
1:B:357:GLN:HE22	1:B:391:THR:H	1.63	0.46
1:B:357:GLN:NE2	1:B:391:THR:OG1	2.49	0.46
1:F:338:VAL:HG22	1:F:373:THR:HG22	1.98	0.46
1:B:208:ASN:HD22	1:B:296:GLY:HA3	1.79	0.46
1:C:172:LEU:HD13	1:C:204:VAL:HG22	1.96	0.46
1:E:265:ILE:HD13	1:E:272:ILE:HG12	1.97	0.46
1:E:296:GLY:O	1:F:115:GLN:N	2.49	0.46
1:F:46:LEU:HD21	1:F:67:LEU:HD21	1.98	0.46
1:A:366:PRO:HG2	6:D:505:MAN:H61	1.97	0.46
1:C:280:PHE:HD2	1:C:308:VAL:HG23	1.80	0.46
1:D:335:VAL:HG22	1:D:375:VAL:HG22	1.97	0.46
1:E:126:LEU:HD11	1:E:166:LEU:HB2	1.96	0.46
1:C:80:ILE:HG13	1:C:89:PHE:HB2	1.98	0.45
1:D:157:ARG:HD3	1:D:158:SER:N	2.31	0.45
1:E:179:GLU:OE2	1:E:201:ARG:HD2	2.15	0.45
1:A:64:ARG:HD2	1:A:68:CYS:HB2	1.98	0.45
1:E:338:VAL:HG22	1:E:373:THR:HG23	1.98	0.45
1:A:208:ASN:HA	1:A:242:ASP:OD2	2.17	0.45
1:C:115:GLN:N	1:D:296:GLY:O	2.44	0.45
1:C:322:SER:O	1:D:37:ARG:NH2	2.50	0.45
1:B:179:GLU:OE2	1:B:201:ARG:HD3	2.17	0.44
1:C:265:ILE:HD13	1:C:272:ILE:HG12	1.99	0.44
1:D:224:PRO:HA	1:D:309:LEU:HB2	1.98	0.44
1:D:3:ILE:HG23	1:E:148:LEU:HD13	1.99	0.44
1:D:77:HIS:O	1:D:77:HIS:ND1	2.51	0.44
6:A:506:MAN:O2	6:A:507:MAN:O4	2.27	0.44
1:E:29:GLY:O	1:E:33:THR:OG1	2.30	0.44
1:E:357:GLN:NE2	1:E:391:THR:OG1	2.50	0.44
1:A:320:ILE:HG23	1:A:339:PHE:HB3	1.99	0.44
1:B:9:GLU:HB3	1:B:98:ILE:HG22	2.00	0.44
1:D:9:GLU:HG2	1:D:10:GLU:HG2	2.00	0.44
1:E:216:GLN:HB2	1:E:219:TYR:CZ	2.53	0.44
1:C:208:ASN:HD22	1:C:296:GLY:HA3	1.82	0.44
1:D:280:PHE:HD2	1:D:308:VAL:HG23	1.83	0.44
1:A:180:LEU:HB2	1:A:202:ILE:HB	1.98	0.44
1:C:192:PRO:HA	1:C:193:PRO:HD3	1.89	0.44
1:A:414:ASP:OD2	1:A:415:ILE:N	2.47	0.44
1:C:64:ARG:HD2	1:C:68:CYS:HB2	2.00	0.44
1:D:387:GLU:HA	1:D:411:GLN:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:ASN:OD1	1:E:142:TYR:N	2.49	0.44
1:E:88:PHE:CE1	1:F:338:VAL:HG21	2.53	0.44
1:E:254:PHE:CE2	1:F:125:PRO:HB2	2.48	0.44
1:D:219:TYR:HB2	1:D:304:VAL:HG22	1.99	0.43
1:C:64:ARG:HH12	1:C:134:ILE:HG23	1.83	0.43
1:F:131:ASP:HB2	1:F:139:VAL:HG23	2.01	0.43
1:A:141:ASN:HB3	1:A:187:LEU:HB2	2.00	0.43
1:B:116:GLU:HG3	1:B:173:ASP:HA	2.00	0.43
1:B:265:ILE:HD13	1:B:272:ILE:HG12	2.01	0.43
1:C:145:SER:HB3	1:C:183:ILE:HB	2.01	0.43
1:C:224:PRO:HA	1:C:309:LEU:HB2	2.00	0.43
1:D:219:TYR:CZ	1:D:236:VAL:HG23	2.54	0.43
1:E:271:GLU:OE1	1:E:273:ARG:NH2	2.52	0.43
1:E:208:ASN:HB2	1:E:246:HIS:ND1	2.34	0.43
1:B:40:HIS:CE1	1:B:45:GLU:HA	2.53	0.42
1:D:208:ASN:HD22	1:D:296:GLY:HA3	1.84	0.42
1:F:188:ASP:HB3	1:F:194:LYS:H	1.83	0.42
1:E:208:ASN:HD22	1:E:296:GLY:HA3	1.83	0.42
1:F:9:GLU:HG2	1:F:10:GLU:HG2	2.00	0.42
1:A:9:GLU:HG2	1:A:10:GLU:HG2	2.00	0.42
1:A:9:GLU:HB3	1:A:98:ILE:HG12	2.01	0.42
1:D:34:ARG:HG2	1:D:84:ASN:O	2.19	0.42
1:E:214:PHE:CD2	1:E:302:CYS:HB3	2.54	0.42
1:F:64:ARG:HH22	1:F:134:ILE:HG23	1.85	0.42
1:A:366:PRO:HB3	1:A:372:TYR:CE1	2.54	0.42
1:C:360:LEU:HA	1:C:361:PRO:HD3	1.80	0.42
1:D:340:SER:HB3	1:D:371:PHE:HD1	1.84	0.42
1:D:318:LEU:HD21	1:D:320:ILE:HD11	2.02	0.42
1:E:30:GLU:O	1:E:34:ARG:HG3	2.20	0.42
1:F:312:ASN:ND2	1:F:400:PRO:O	2.52	0.42
1:B:9:GLU:HG2	1:B:10:GLU:N	2.34	0.42
1:C:254:PHE:CE2	1:D:125:PRO:HB2	2.55	0.42
1:D:65:GLU:HG2	1:D:132:PRO:HB2	2.01	0.42
1:D:141:ASN:HB3	1:D:187:LEU:HB2	2.02	0.42
1:D:188:ASP:HB3	1:D:194:LYS:H	1.85	0.42
1:B:112:LEU:HB2	1:B:202:ILE:HG12	2.01	0.42
1:C:191:ALA:HA	1:C:192:PRO:HA	1.88	0.42
1:C:258:GLY:HA2	1:D:130:GLN:HG2	2.02	0.42
1:D:145:SER:HA	1:D:146:PRO:HD3	1.92	0.42
1:B:208:ASN:HB2	1:B:246:HIS:ND1	2.34	0.42
1:E:149:HIS:O	1:E:168:LEU:HD12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:340:SER:HB3	1:C:371:PHE:CD1	2.54	0.42
1:B:64:ARG:HD2	1:B:68:CYS:HB2	2.02	0.41
1:E:191:ALA:HA	1:E:192:PRO:HA	1.87	0.41
1:E:214:PHE:CG	1:E:302:CYS:HB3	2.55	0.41
1:A:360:LEU:HA	1:A:361:PRO:HD3	1.79	0.41
1:B:262:PRO:HB2	1:B:276:GLY:H	1.84	0.41
1:B:387:GLU:HB3	1:B:411:GLN:HG2	2.03	0.41
1:C:208:ASN:HB2	1:C:246:HIS:ND1	2.35	0.41
1:C:317:LYS:HB2	1:C:342:SER:HB3	2.01	0.41
1:C:254:PHE:HE2	1:D:125:PRO:HB2	1.85	0.41
1:D:7:MET:HA	1:D:8:PRO:HD3	1.95	0.41
1:E:227:SER:HA	1:E:228:PRO:HD3	1.92	0.41
1:A:145:SER:HB3	1:A:183:ILE:HB	2.02	0.41
1:B:210:ASN:ND2	1:B:242:ASP:OD1	2.50	0.41
1:D:84:ASN:HA	1:D:85:PRO:HA	1.81	0.41
1:E:131:ASP:HB2	1:E:139:VAL:HG13	2.03	0.41
1:E:135:GLY:O	1:E:140:GLN:NE2	2.53	0.41
1:F:15:TYR:HB3	1:F:58:LEU:HD12	2.02	0.41
1:F:7:MET:HA	1:F:8:PRO:HD3	1.90	0.41
1:B:286:TYR:HB2	1:B:306:ILE:HG23	2.01	0.41
1:B:117:ILE:HD13	1:B:207:ILE:HG21	2.03	0.41
1:B:208:ASN:ND2	1:B:296:GLY:HA3	2.35	0.41
1:F:116:GLU:HG3	1:F:173:ASP:HA	2.03	0.41
1:A:79:GLN:HG3	1:B:371:PHE:CE1	2.56	0.41
1:A:7:MET:HA	1:A:8:PRO:HD3	1.94	0.41
1:C:34:ARG:O	1:C:83:GLU:HG2	2.20	0.41
1:E:294:ASP:OD2	1:E:298:LEU:HB2	2.21	0.41
1:F:374:LEU:HD21	1:F:390:ILE:HG21	2.01	0.41
1:C:9:GLU:HG2	1:C:10:GLU:HG2	2.03	0.41
1:D:63:ASP:OD2	1:D:66:ALA:N	2.52	0.41
1:E:290:ILE:HB	1:E:302:CYS:SG	2.61	0.41
1:E:40:HIS:CE1	1:E:45:GLU:HG2	2.56	0.41
1:E:9:GLU:HG2	1:E:10:GLU:HG2	2.03	0.41
1:E:211:ALA:HA	1:E:212:PRO:HD3	1.96	0.41
1:A:105:PHE:HA	1:A:106:PRO:HD3	1.96	0.40
1:D:76:LEU:HA	1:D:76:LEU:HD12	1.94	0.40
1:A:208:ASN:HB3	1:B:115:GLN:OE1	2.21	0.40
1:C:227:SER:HA	1:C:228:PRO:HD3	1.95	0.40
1:E:378:ARG:HH12	1:E:380:LEU:HD23	1.87	0.40
1:F:279:ASP:OD2	1:F:282:ALA:N	2.52	0.40
1:B:82:LEU:O	1:B:86:VAL:HA	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:208:ASN:HB2	1:F:246:HIS:ND1	2.36	0.40
1:A:145:SER:HA	1:A:146:PRO:HD3	1.91	0.40
1:A:371:PHE:CZ	1:B:79:GLN:HG3	2.56	0.40
1:B:112:LEU:HD11	1:B:126:LEU:HD21	2.01	0.40
1:B:280:PHE:HD1	1:B:308:VAL:HG23	1.86	0.40
1:D:123:VAL:HG12	1:D:167:VAL:HG22	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	412/426 (97%)	389 (94%)	23 (6%)	0	100	100
1	B	413/426 (97%)	395 (96%)	18 (4%)	0	100	100
1	C	413/426 (97%)	393 (95%)	20 (5%)	0	100	100
1	D	410/426 (96%)	388 (95%)	22 (5%)	0	100	100
1	E	411/426 (96%)	391 (95%)	20 (5%)	0	100	100
1	F	413/426 (97%)	388 (94%)	25 (6%)	0	100	100
All	All	2472/2556 (97%)	2344 (95%)	128 (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	339/370 (92%)	337 (99%)	2 (1%)	86	96
1	B	337/370 (91%)	335 (99%)	2 (1%)	86	96
1	C	330/370 (89%)	328 (99%)	2 (1%)	86	96
1	D	330/370 (89%)	329 (100%)	1 (0%)	92	98
1	E	341/370 (92%)	341 (100%)	0	100	100
1	F	345/370 (93%)	344 (100%)	1 (0%)	92	98
All	All	2022/2220 (91%)	2014 (100%)	8 (0%)	91	97

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	148	LEU
1	B	157	ARG
1	B	288	MET
1	C	159	ASP
1	C	162	LYS
1	D	157	ARG
1	F	200	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	149	HIS

### 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 ⓘ

23 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	G	1	1,2	14,14,15	0.24	0	17,19,21	0.37	0
2	FUC	G	2	2	10,10,11	0.74	0	14,14,16	0.94	0
2	NAG	H	1	1,2	14,14,15	0.22	0	17,19,21	0.43	0
2	FUC	H	2	2	10,10,11	0.77	0	14,14,16	0.82	0
3	NAG	I	1	1,3	14,14,15	0.41	0	17,19,21	0.42	0
3	NAG	I	2	3	14,14,15	0.22	0	17,19,21	0.39	0
3	BMA	I	3	3	11,11,12	1.03	0	15,15,17	1.68	3 (20%)
3	MAN	I	4	3	11,11,12	0.83	1 (9%)	15,15,17	1.39	2 (13%)
3	NAG	I	5	3	14,14,15	0.28	0	17,19,21	0.49	0
3	GAL	I	6	3	11,11,12	0.55	0	15,15,17	0.90	0
3	MAN	I	7	3	11,11,12	1.31	2 (18%)	15,15,17	1.12	2 (13%)
3	NAG	I	8	3	14,14,15	0.21	0	17,19,21	0.45	0
3	FUC	I	9	3	10,10,11	0.72	0	14,14,16	0.85	0
4	NAG	J	1	1,4	14,14,15	0.40	0	17,19,21	0.65	0
4	NAG	J	2	4	14,14,15	0.58	1 (7%)	17,19,21	0.53	0
4	BMA	J	3	4	11,11,12	0.73	0	15,15,17	0.73	0
2	NAG	K	1	1,2	14,14,15	0.31	0	17,19,21	0.38	0
2	FUC	K	2	2	10,10,11	0.78	0	14,14,16	0.80	0
5	NAG	L	1	1,5	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	L	2	5	14,14,15	0.27	0	17,19,21	0.41	0
5	BMA	L	3	5	11,11,12	0.61	0	15,15,17	0.91	0
5	MAN	L	4	5	11,11,12	0.65	0	15,15,17	1.12	2 (13%)
5	FUC	L	5	5	10,10,11	0.73	0	14,14,16	0.84	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	G	1	1,2	-	0/6/23/26	0/1/1/1
2	FUC	G	2	2	-	-	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	FUC	H	2	2	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	I	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1
3	BMA	I	3	3	-	2/2/19/22	1/1/1/1
3	MAN	I	4	3	-	1/2/19/22	0/1/1/1
3	NAG	I	5	3	-	0/6/23/26	0/1/1/1
3	GAL	I	6	3	-	2/2/19/22	0/1/1/1
3	MAN	I	7	3	-	2/2/19/22	0/1/1/1
3	NAG	I	8	3	-	0/6/23/26	0/1/1/1
3	FUC	I	9	3	-	-	0/1/1/1
4	NAG	J	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	2/2/19/22	0/1/1/1
2	NAG	K	1	1,2	-	0/6/23/26	0/1/1/1
2	FUC	K	2	2	-	-	0/1/1/1
5	NAG	L	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	L	2	5	-	2/6/23/26	0/1/1/1
5	BMA	L	3	5	-	1/2/19/22	0/1/1/1
5	MAN	L	4	5	-	1/2/19/22	0/1/1/1
5	FUC	L	5	5	-	-	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	7	MAN	O5-C1	-3.18	1.38	1.43
3	I	4	MAN	O5-C5	2.31	1.48	1.43
3	I	7	MAN	C4-C3	2.24	1.58	1.52
4	J	2	NAG	O5-C1	-2.04	1.40	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	4	MAN	C1-O5-C5	4.44	118.21	112.19
3	I	3	BMA	C1-O5-C5	4.16	117.83	112.19
3	I	3	BMA	C1-C2-C3	3.68	114.19	109.67
5	L	4	MAN	C1-O5-C5	2.64	115.78	112.19
3	I	3	BMA	O2-C2-C3	-2.50	105.12	110.14
3	I	4	MAN	O2-C2-C3	-2.25	105.62	110.14
3	I	7	MAN	O2-C2-C3	-2.24	105.65	110.14
5	L	4	MAN	O2-C2-C3	-2.24	105.65	110.14
3	I	7	MAN	C2-C3-C4	2.04	114.43	110.89

There are no chirality outliers.

All (16) torsion outliers are listed below:

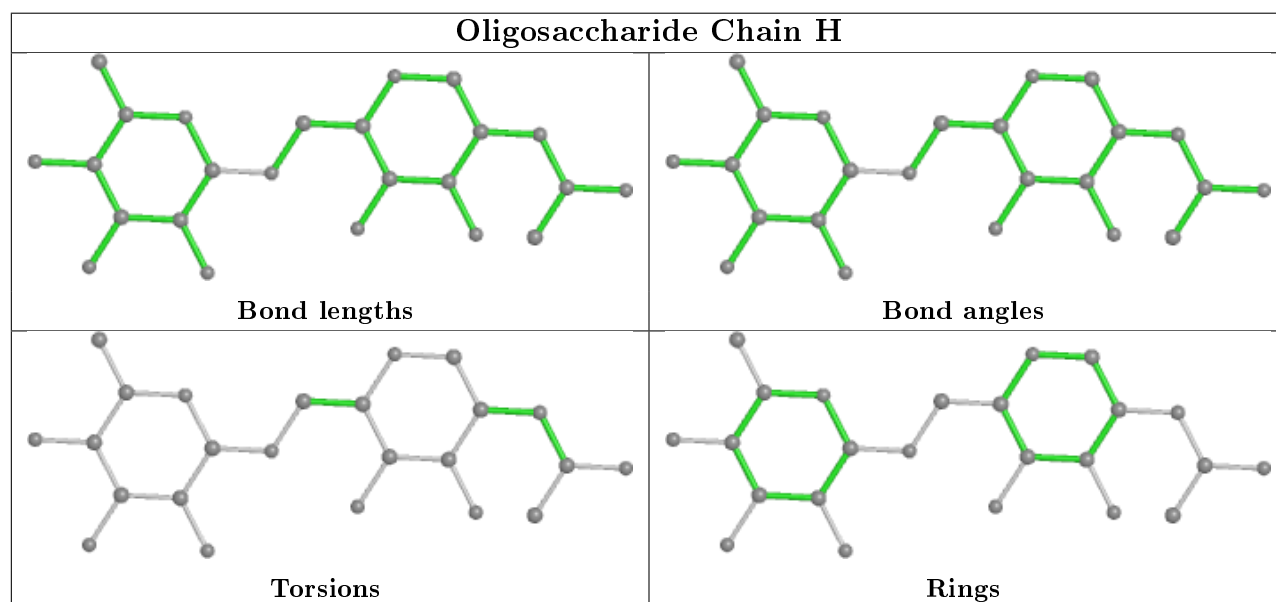
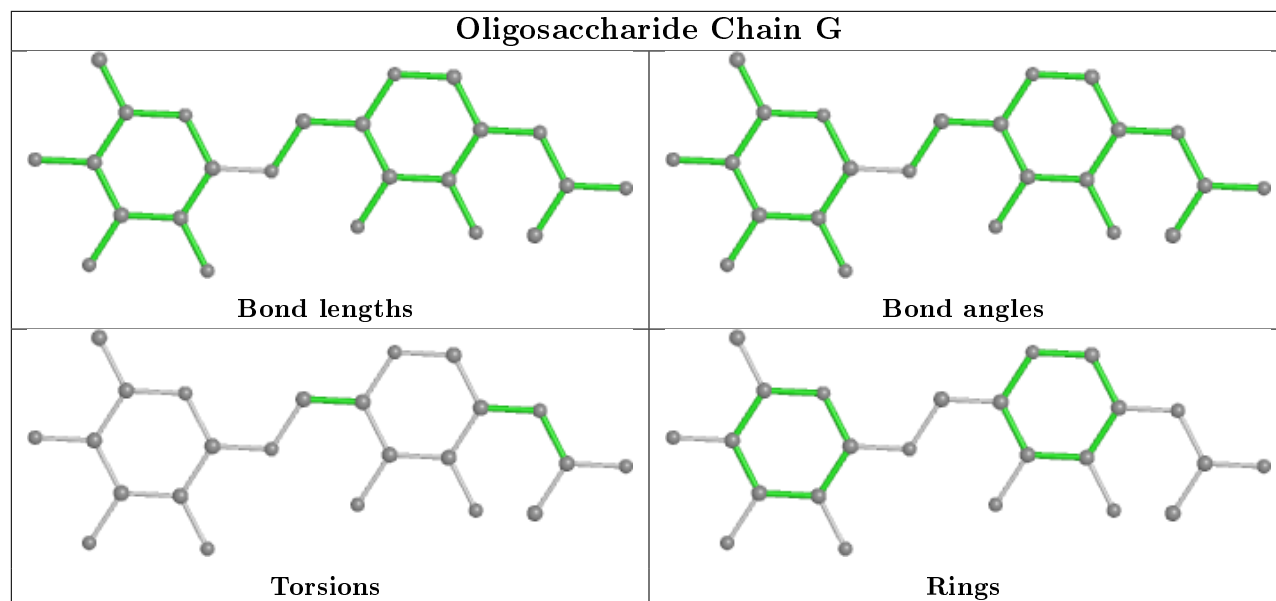
Mol	Chain	Res	Type	Atoms
5	L	2	NAG	O5-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6
5	L	2	NAG	C4-C5-C6-O6
4	J	3	BMA	O5-C5-C6-O6
3	I	3	BMA	O5-C5-C6-O6
3	I	2	NAG	C4-C5-C6-O6
5	L	4	MAN	O5-C5-C6-O6
4	J	1	NAG	O5-C5-C6-O6
3	I	4	MAN	O5-C5-C6-O6
3	I	3	BMA	C4-C5-C6-O6
3	I	7	MAN	O5-C5-C6-O6
3	I	7	MAN	C4-C5-C6-O6
3	I	6	GAL	O5-C5-C6-O6
4	J	3	BMA	C4-C5-C6-O6
5	L	3	BMA	O5-C5-C6-O6
3	I	6	GAL	C4-C5-C6-O6

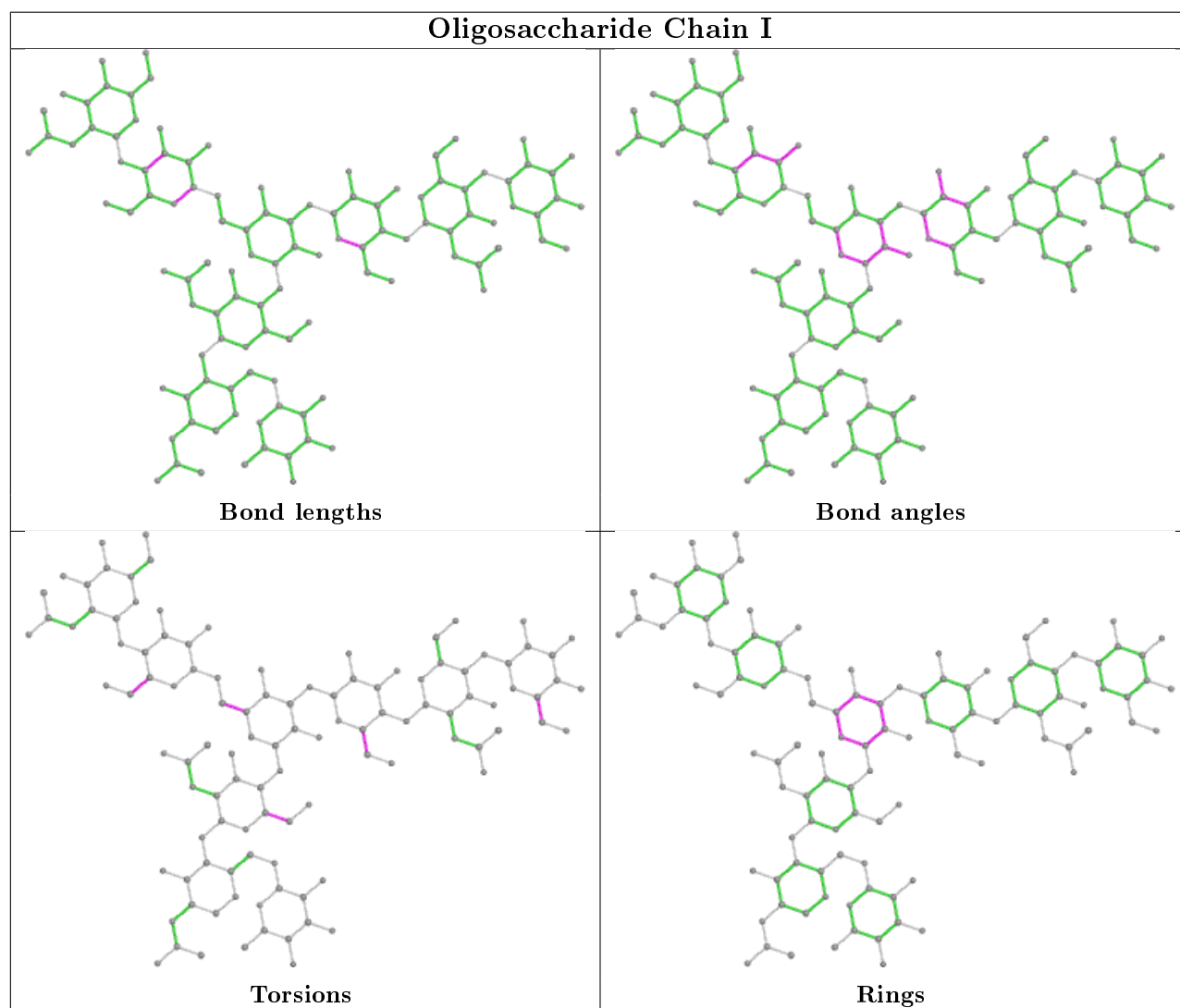
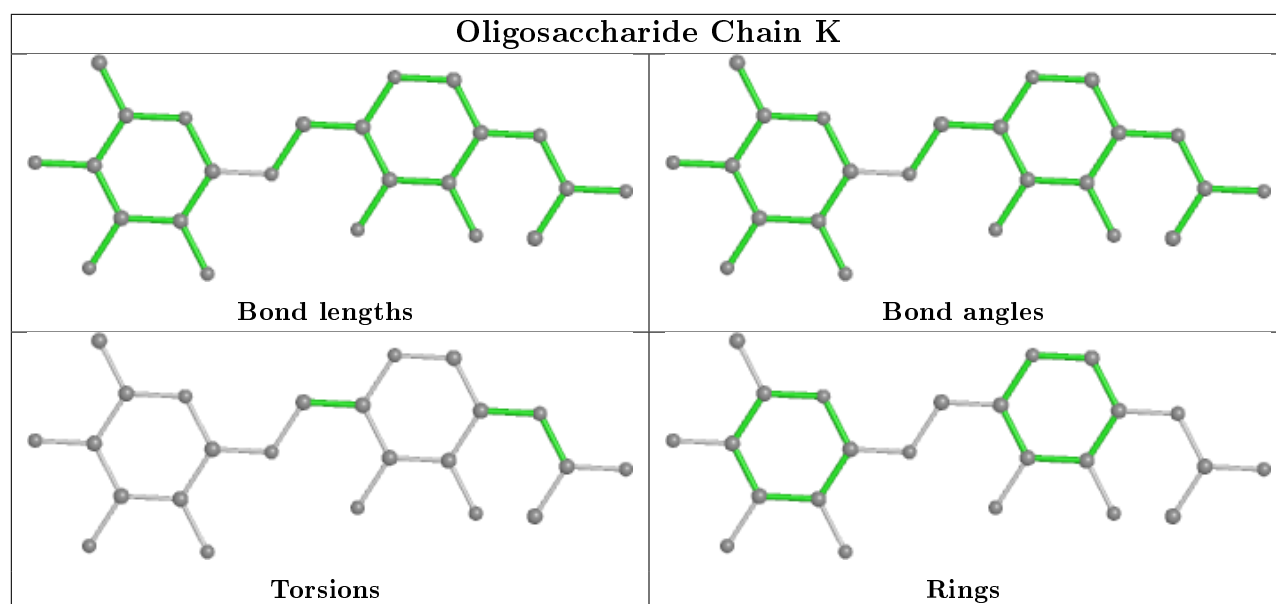
All (1) ring outliers are listed below:

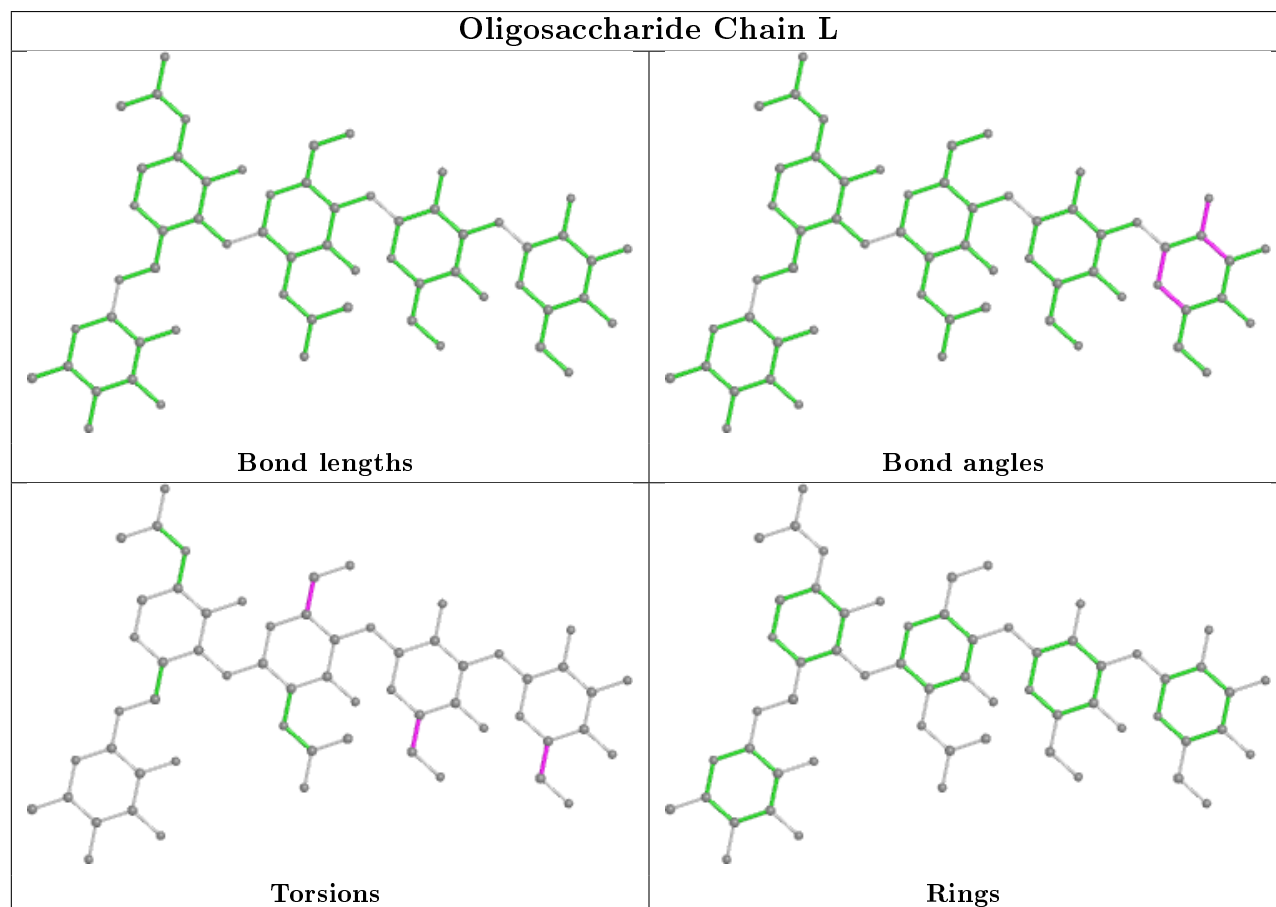
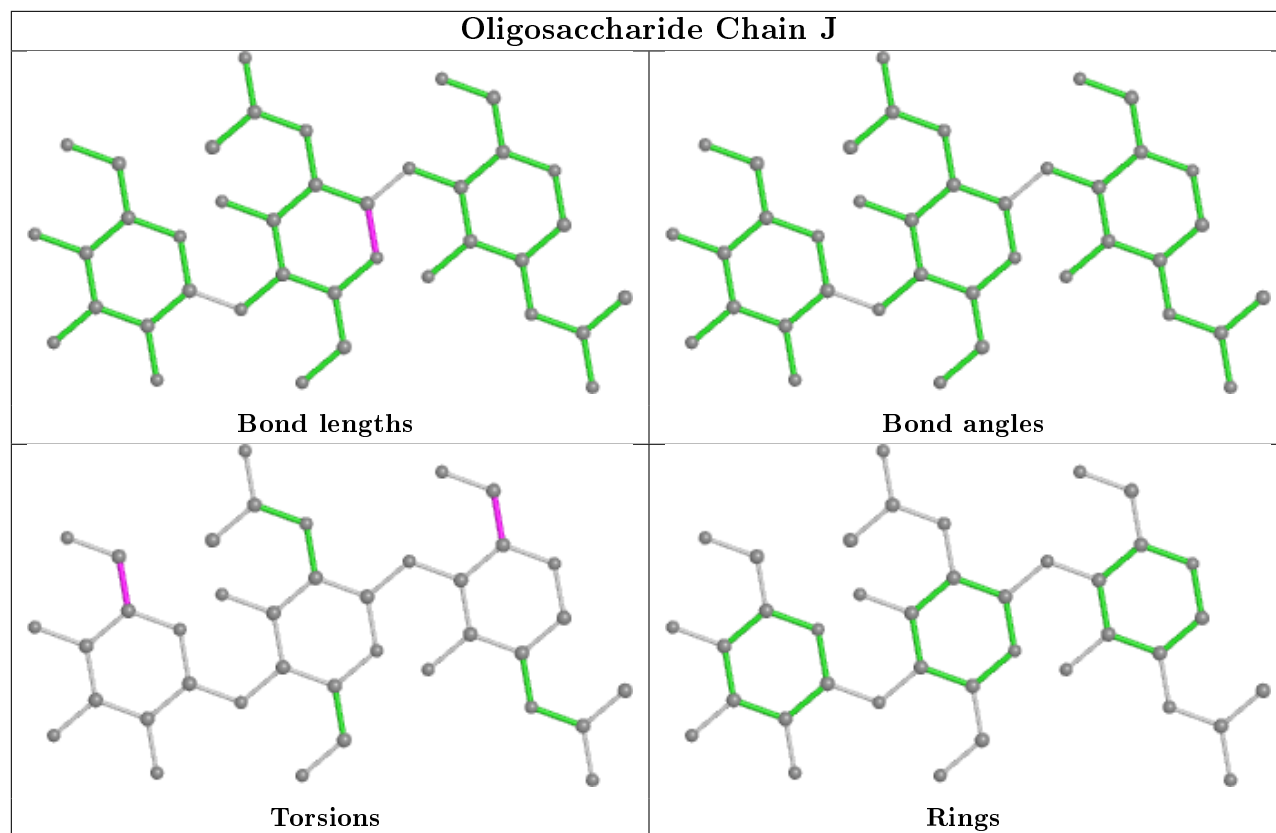
Mol	Chain	Res	Type	Atoms
3	I	3	BMA	C1-C2-C3-C4-C5-O5

No monomer is involved in short contacts.

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 75 ligands modelled in this entry, 54 are monoatomic - leaving 21 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$
8	NAG	F	502	1	14,14,15	0.25	0	17,19,21	0.48	0
6	MAN	D	505	1	11,11,12	0.73	0	15,15,17	0.99	1 (6%)
6	MAN	F	505	1	11,11,12	0.73	0	15,15,17	1.02	2 (13%)
6	MAN	E	507	1	11,11,12	0.72	0	15,15,17	1.07	2 (13%)
6	MAN	C	506	1	11,11,12	0.80	0	15,15,17	0.93	2 (13%)
6	MAN	B	511	1	11,11,12	0.73	0	15,15,17	1.01	2 (13%)
6	MAN	A	506	1	11,11,12	0.78	0	15,15,17	1.06	2 (13%)
6	MAN	E	506	1	11,11,12	0.61	0	15,15,17	1.34	3 (20%)
6	MAN	F	503	1	11,11,12	0.83	1 (9%)	15,15,17	1.72	3 (20%)
6	MAN	A	505	1	11,11,12	0.60	0	15,15,17	1.22	2 (13%)
6	MAN	D	503	1	11,11,12	1.10	0	15,15,17	1.30	2 (13%)
6	MAN	C	505	1	11,11,12	0.79	0	15,15,17	1.19	3 (20%)
6	MAN	E	508	1	11,11,12	0.79	0	15,15,17	0.94	1 (6%)
8	NAG	B	501	1	14,14,15	0.22	0	17,19,21	0.37	0
6	MAN	C	504	1	11,11,12	0.62	0	15,15,17	1.21	2 (13%)
6	MAN	B	513	1	11,11,12	0.84	0	15,15,17	0.90	1 (6%)
8	NAG	F	501	1	14,14,15	0.18	0	17,19,21	0.41	0
6	MAN	D	504	1	11,11,12	0.65	0	15,15,17	1.16	2 (13%)
6	MAN	F	504	1	11,11,12	0.72	0	15,15,17	1.09	2 (13%)
6	MAN	A	507	1	11,11,12	0.79	0	15,15,17	1.03	1 (6%)
6	MAN	B	512	1	11,11,12	0.74	0	15,15,17	1.25	3 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	F	502	1	-	2/6/23/26	0/1/1/1
6	MAN	D	505	1	-	0/2/19/22	0/1/1/1
6	MAN	F	505	1	-	0/2/19/22	0/1/1/1
6	MAN	E	507	1	-	0/2/19/22	0/1/1/1
6	MAN	C	506	1	-	0/2/19/22	0/1/1/1
6	MAN	B	511	1	-	0/2/19/22	0/1/1/1
6	MAN	A	506	1	-	2/2/19/22	0/1/1/1
6	MAN	E	506	1	-	0/2/19/22	0/1/1/1
6	MAN	F	503	1	-	0/2/19/22	0/1/1/1
6	MAN	A	505	1	-	0/2/19/22	0/1/1/1
6	MAN	D	503	1	-	0/2/19/22	0/1/1/1
6	MAN	C	505	1	-	0/2/19/22	0/1/1/1
6	MAN	E	508	1	-	1/2/19/22	0/1/1/1
8	NAG	B	501	1	-	0/6/23/26	0/1/1/1
6	MAN	C	504	1	-	0/2/19/22	0/1/1/1
6	MAN	B	513	1	-	2/2/19/22	0/1/1/1
8	NAG	F	501	1	-	3/6/23/26	0/1/1/1
6	MAN	D	504	1	-	0/2/19/22	0/1/1/1
6	MAN	F	504	1	-	0/2/19/22	0/1/1/1
6	MAN	A	507	1	-	1/2/19/22	0/1/1/1
6	MAN	B	512	1	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	503	MAN	O5-C5	2.25	1.48	1.43

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	503	MAN	C1-O5-C5	5.19	119.22	112.19
6	E	506	MAN	C1-O5-C5	3.63	117.11	112.19
6	C	504	MAN	C1-O5-C5	3.22	116.56	112.19
6	A	505	MAN	C1-O5-C5	3.18	116.51	112.19
6	D	504	MAN	C1-O5-C5	2.86	116.07	112.19
6	D	503	MAN	C1-O5-C5	2.69	115.83	112.19
6	F	503	MAN	O5-C1-C2	2.63	114.83	110.77
6	B	512	MAN	C1-O5-C5	2.63	115.75	112.19
6	B	512	MAN	O5-C1-C2	2.39	114.46	110.77
6	F	504	MAN	C1-O5-C5	2.38	115.42	112.19
6	C	505	MAN	C1-O5-C5	2.34	115.36	112.19
6	F	505	MAN	C1-O5-C5	2.34	115.36	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	E	506	MAN	O5-C1-C2	2.30	114.31	110.77
6	E	507	MAN	C1-O5-C5	2.27	115.27	112.19
6	E	507	MAN	O2-C2-C3	-2.27	105.59	110.14
6	B	511	MAN	C1-O5-C5	2.25	115.25	112.19
6	A	505	MAN	O2-C2-C3	-2.25	105.63	110.14
6	D	505	MAN	O2-C2-C3	-2.25	105.63	110.14
6	B	513	MAN	O2-C2-C3	-2.24	105.65	110.14
6	E	506	MAN	O2-C2-C3	-2.23	105.66	110.14
6	C	506	MAN	O2-C2-C3	-2.23	105.67	110.14
6	C	504	MAN	O2-C2-C3	-2.23	105.67	110.14
6	A	506	MAN	O2-C2-C3	-2.22	105.68	110.14
6	F	504	MAN	O2-C2-C3	-2.19	105.74	110.14
6	E	508	MAN	O2-C2-C3	-2.19	105.75	110.14
6	A	507	MAN	O2-C2-C3	-2.19	105.75	110.14
6	C	505	MAN	O5-C1-C2	2.18	114.14	110.77
6	F	505	MAN	O2-C2-C3	-2.18	105.77	110.14
6	B	511	MAN	O2-C2-C3	-2.17	105.79	110.14
6	D	504	MAN	O2-C2-C3	-2.16	105.82	110.14
6	C	506	MAN	C1-O5-C5	2.14	115.09	112.19
6	F	503	MAN	O2-C2-C3	-2.10	105.94	110.14
6	B	512	MAN	O2-C2-C3	-2.09	105.94	110.14
6	C	505	MAN	O2-C2-C3	-2.08	105.97	110.14
6	A	506	MAN	C1-O5-C5	2.06	114.99	112.19
6	D	503	MAN	O2-C2-C3	-2.01	106.11	110.14

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	F	502	NAG	O5-C5-C6-O6
6	B	513	MAN	C4-C5-C6-O6
8	F	501	NAG	C8-C7-N2-C2
8	F	501	NAG	O7-C7-N2-C2
8	F	502	NAG	C4-C5-C6-O6
6	A	506	MAN	O5-C5-C6-O6
6	B	513	MAN	O5-C5-C6-O6
6	E	508	MAN	O5-C5-C6-O6
6	A	506	MAN	C4-C5-C6-O6
8	F	501	NAG	O5-C5-C6-O6
6	A	507	MAN	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	505	MAN	1	0
6	A	506	MAN	2	0
6	A	507	MAN	2	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, 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	414/426 (97%)	0.04	13 (3%) 49 44	48, 108, 171, 229	0
1	B	415/426 (97%)	0.14	23 (5%) 25 21	53, 103, 165, 225	0
1	C	415/426 (97%)	0.40	44 (10%) 6 4	53, 114, 219, 343	0
1	D	412/426 (96%)	0.29	35 (8%) 10 8	52, 137, 223, 383	0
1	E	413/426 (96%)	0.02	6 (1%) 73 73	53, 108, 170, 267	0
1	F	415/426 (97%)	0.14	23 (5%) 25 21	58, 98, 158, 247	0
All	All	2484/2556 (97%)	0.17	144 (5%) 23 19	48, 108, 192, 383	0

All (144) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	18	ALA	9.3
1	C	331	ALA	8.8
1	C	337	ALA	8.6
1	D	82	LEU	7.5
1	D	337	ALA	7.4
1	D	81	ILE	6.1
1	C	81	ILE	6.1
1	C	338	VAL	5.9
1	C	332	PRO	5.7
1	F	2	ALA	5.5
1	D	78	PHE	5.3
1	D	320	ILE	5.3
1	C	80	ILE	5.2
1	D	2	ALA	4.9
1	A	88	PHE	4.8
1	C	82	LEU	4.8
1	C	333	GLU	4.7
1	F	3	ILE	4.6
1	D	89	PHE	4.6

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Mol	Chain	Res	Type	RSRZ
1	D	40	HIS	4.5
1	D	19	ASN	4.5
1	D	11	THR	4.4
1	C	375	VAL	4.4
1	D	15	TYR	4.4
1	C	392	ILE	4.3
1	C	95	PHE	4.3
1	B	258	GLY	4.2
1	C	75	VAL	4.2
1	C	393	THR	4.1
1	D	330	ASN	4.1
1	D	5	TYR	4.1
1	F	220	ALA	4.1
1	B	229	LEU	4.1
1	B	230	ASN	4.0
1	C	335	VAL	3.9
1	A	3	ILE	3.9
1	C	339	PHE	3.8
1	B	82	LEU	3.7
1	C	374	LEU	3.7
1	D	80	ILE	3.6
1	B	390	ILE	3.5
1	C	390	ILE	3.5
1	A	80	ILE	3.5
1	F	306	ILE	3.4
1	F	336	VAL	3.4
1	D	36	ALA	3.3
1	D	3	ILE	3.3
1	D	371	PHE	3.2
1	D	390	ILE	3.2
1	D	46	LEU	3.2
1	B	221	VAL	3.2
1	B	320	ILE	3.2
1	F	219	TYR	3.1
1	A	47	LEU	3.1
1	C	334	ALA	3.0
1	B	220	ALA	3.0
1	C	340	SER	3.0
1	F	257	GLY	3.0
1	E	80	ILE	3.0
1	A	82	LEU	3.0
1	B	276	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	93	LEU	3.0
1	B	257	GLY	2.9
1	A	187	LEU	2.9
1	B	361	PRO	2.9
1	F	276	GLY	2.9
1	B	89	PHE	2.8
1	D	374	LEU	2.8
1	F	256	GLY	2.8
1	F	172	LEU	2.8
1	F	78	PHE	2.8
1	C	47	LEU	2.7
1	B	363	LEU	2.7
1	D	17	VAL	2.7
1	C	415	ILE	2.7
1	A	86	VAL	2.7
1	C	371	PHE	2.6
1	D	150	PHE	2.6
1	C	318	LEU	2.6
1	D	406	HIS	2.6
1	C	24	LEU	2.6
1	C	323	LEU	2.6
1	F	360	LEU	2.5
1	C	230	ASN	2.5
1	D	96	THR	2.5
1	C	408	ILE	2.5
1	F	339	PHE	2.4
1	E	38	ILE	2.4
1	E	172	LEU	2.4
1	F	324	THR	2.4
1	D	338	VAL	2.4
1	D	182	LEU	2.4
1	B	339	PHE	2.4
1	C	223	VAL	2.4
1	A	356	ILE	2.4
1	B	272	ILE	2.4
1	C	276	GLY	2.4
1	C	355	SER	2.4
1	F	327	ILE	2.4
1	F	390	ILE	2.4
1	C	364	LEU	2.3
1	D	16	LEU	2.3
1	C	336	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	241	LEU	2.3
1	D	79	GLN	2.3
1	A	172	LEU	2.3
1	A	320	ILE	2.3
1	A	17	VAL	2.3
1	A	397	LEU	2.3
1	C	46	LEU	2.3
1	C	325	SER	2.3
1	B	374	LEU	2.2
1	C	93	LEU	2.2
1	C	94	GLN	2.2
1	E	31	LEU	2.2
1	D	75	VAL	2.2
1	D	172	LEU	2.2
1	F	218	LEU	2.2
1	D	86	VAL	2.2
1	D	408	ILE	2.1
1	B	31	LEU	2.1
1	C	397	LEU	2.1
1	A	196	GLY	2.1
1	B	309	LEU	2.1
1	F	337	ALA	2.1
1	C	36	ALA	2.1
1	C	272	ILE	2.1
1	C	410	VAL	2.1
1	E	155	LEU	2.1
1	B	5	TYR	2.1
1	F	148	LEU	2.1
1	F	215	LEU	2.1
1	C	233	VAL	2.1
1	B	233	VAL	2.1
1	E	364	LEU	2.0
1	C	328	PRO	2.0
1	B	323	LEU	2.0
1	B	364	LEU	2.0
1	F	47	LEU	2.0
1	C	117	ILE	2.0
1	F	392	ILE	2.0
1	D	73	PRO	2.0
1	B	305	ALA	2.0
1	C	394	VAL	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 ⓘ

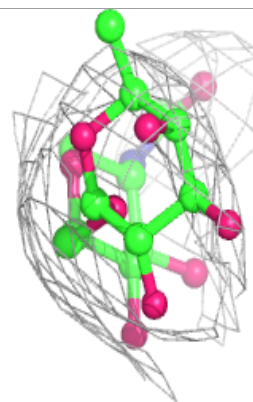
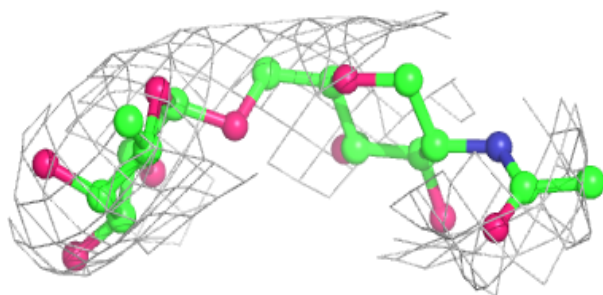
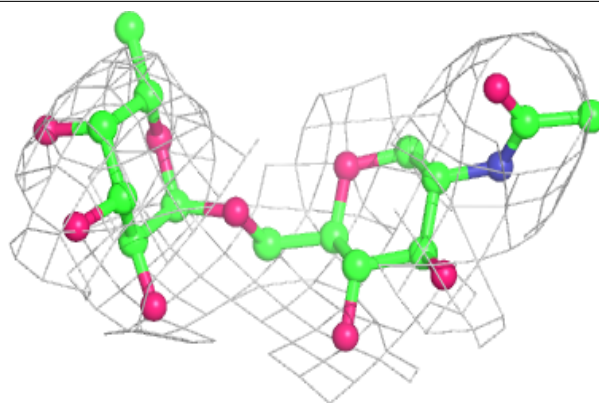
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	BMA	J	3	11/12	0.73	0.18	152,168,177,180	0
2	FUC	H	2	10/11	0.77	0.23	126,145,154,158	0
4	NAG	J	1	14/15	0.79	0.15	108,131,153,161	0
4	NAG	J	2	14/15	0.81	0.20	155,170,174,175	0
3	MAN	I	4	11/12	0.81	0.19	121,154,160,164	0
2	FUC	K	2	10/11	0.82	0.15	145,159,162,163	0
5	BMA	L	3	11/12	0.82	0.10	135,152,155,156	0
3	MAN	I	7	11/12	0.83	0.13	101,134,149,152	0
3	GAL	I	6	11/12	0.86	0.26	141,148,151,151	0
3	BMA	I	3	11/12	0.86	0.08	133,137,154,156	0
2	FUC	G	2	10/11	0.88	0.15	142,149,153,155	0
5	FUC	L	5	10/11	0.89	0.22	140,144,147,151	0
5	NAG	L	2	14/15	0.90	0.13	132,146,152,154	0
2	NAG	G	1	14/15	0.91	0.09	95,130,132,141	0
3	NAG	I	5	14/15	0.91	0.24	137,160,168,169	0
2	NAG	H	1	14/15	0.91	0.09	112,136,143,146	0
5	MAN	L	4	11/12	0.91	0.09	108,140,145,149	0
5	NAG	L	1	14/15	0.92	0.09	87,123,136,139	0
2	NAG	K	1	14/15	0.92	0.11	107,129,138,139	0
3	NAG	I	1	14/15	0.93	0.17	99,114,145,148	0
3	NAG	I	8	14/15	0.94	0.21	129,140,150,151	0
3	NAG	I	2	14/15	0.95	0.13	98,117,124,133	0
3	FUC	I	9	10/11	0.96	0.12	126,133,142,151	0

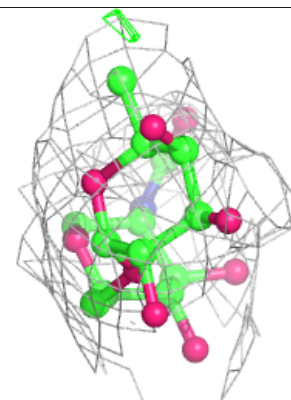
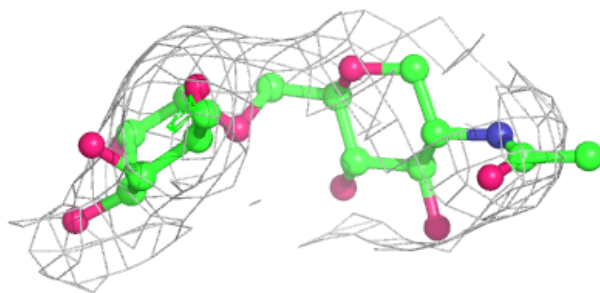
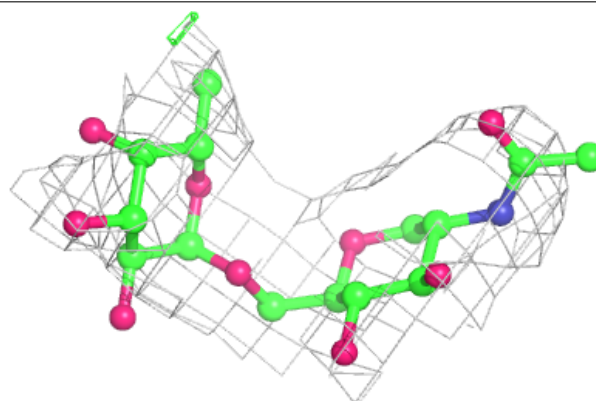
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around Chain G:**

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

**Electron density around Chain H:**

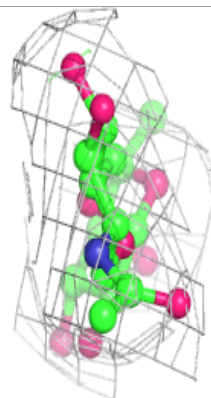
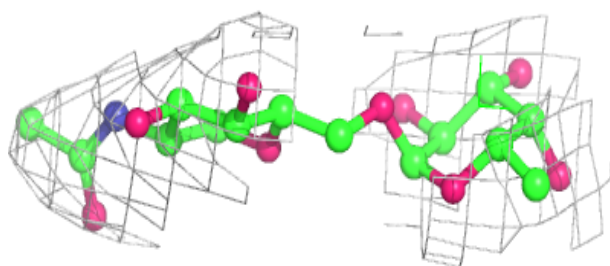
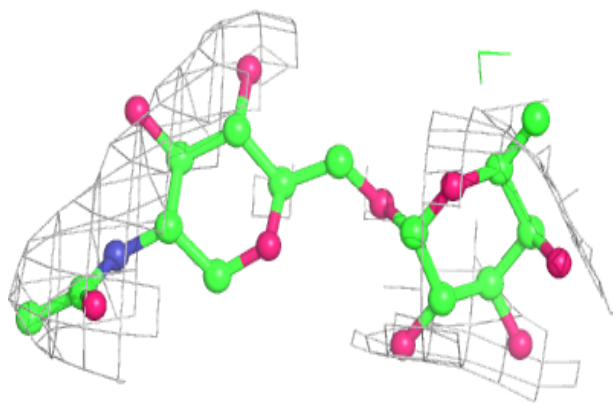
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





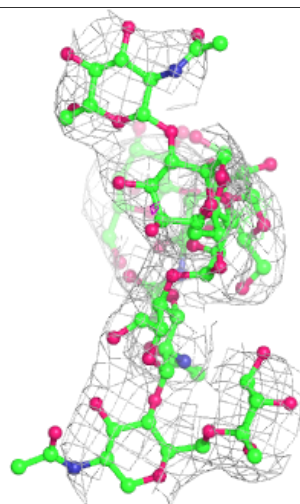
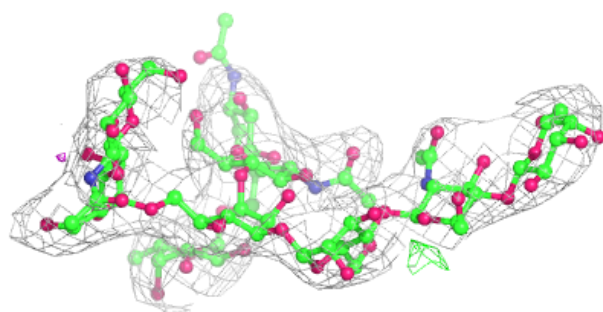
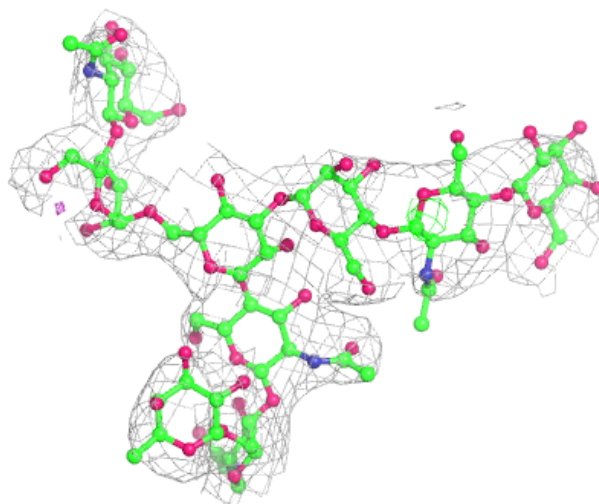
**Electron density around Chain K:**

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



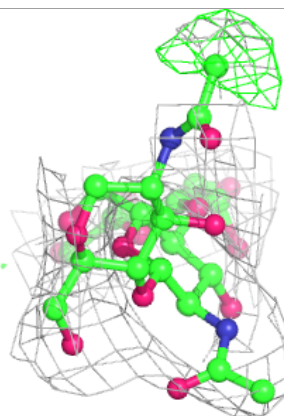
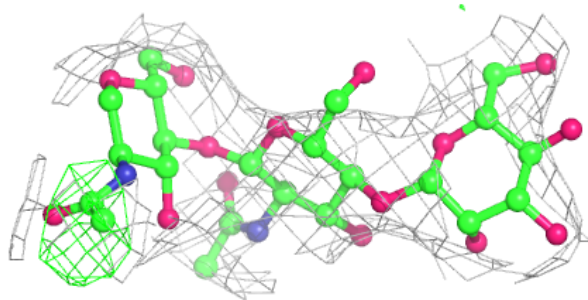
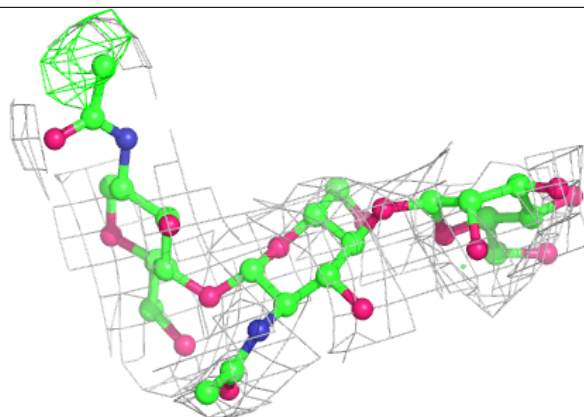
**Electron density around Chain I:**

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

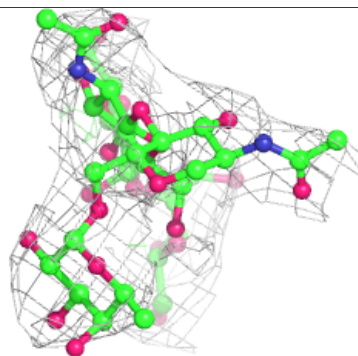
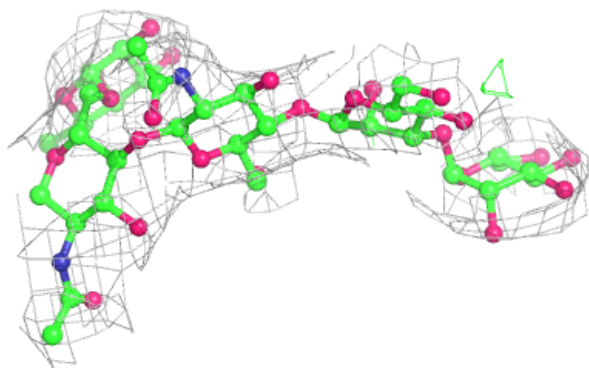
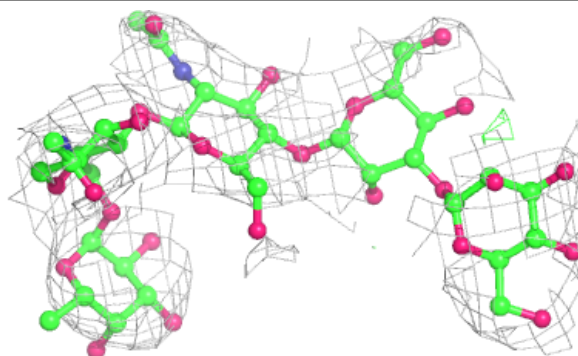


**Electron density around Chain J:**

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

**Electron density around Chain L:**

$2mF_o-DF_c$  (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
8	NAG	F	502	14/15	0.75	0.15	106,131,136,137	0
6	MAN	E	506	11/12	0.80	0.24	97,113,138,139	0
6	MAN	D	503	11/12	0.84	0.18	105,121,132,142	0
7	CA	D	508	1/1	0.84	0.21	119,119,119,119	0
6	MAN	C	506	11/12	0.85	0.20	92,101,118,125	0
6	MAN	A	506	11/12	0.86	0.31	114,136,143,144	0
6	MAN	F	503	11/12	0.86	0.18	93,113,122,125	0
7	CA	D	510	1/1	0.87	0.29	126,126,126,126	0
7	CA	B	517	1/1	0.88	0.38	156,156,156,156	0
7	CA	A	514	1/1	0.88	0.23	52,52,52,52	0
6	MAN	A	507	11/12	0.90	0.21	118,133,139,142	0
6	MAN	C	504	11/12	0.90	0.20	85,98,120,124	0
6	MAN	D	505	11/12	0.92	0.18	110,118,122,124	0
6	MAN	F	505	11/12	0.92	0.13	96,101,112,115	0
8	NAG	F	501	14/15	0.92	0.11	107,126,133,139	0
6	MAN	A	505	11/12	0.92	0.15	105,119,130,144	0
6	MAN	B	511	11/12	0.93	0.22	73,105,108,116	0
7	CA	D	507	1/1	0.93	0.22	89,89,89,89	0
7	CA	E	509	1/1	0.93	0.24	95,95,95,95	0
8	NAG	B	501	14/15	0.94	0.09	87,100,118,118	0
7	CA	A	509	1/1	0.94	0.20	116,116,116,116	0
7	CA	D	506	1/1	0.94	0.17	162,162,162,162	0
7	CA	A	512	1/1	0.94	0.27	53,53,53,53	0
6	MAN	B	513	11/12	0.94	0.16	65,96,123,131	0
7	CA	D	512	1/1	0.95	0.28	81,81,81,81	0
6	MAN	B	512	11/12	0.95	0.15	45,66,91,91	0
7	CA	A	511	1/1	0.95	0.22	83,83,83,83	0
6	MAN	D	504	11/12	0.95	0.16	90,106,122,124	0
7	CA	E	510	1/1	0.95	0.20	66,66,66,66	0
7	CA	B	515	1/1	0.95	0.26	60,60,60,60	0
6	MAN	E	508	11/12	0.95	0.09	83,112,126,129	0
7	CA	B	520	1/1	0.96	0.19	105,105,105,105	0
7	CA	C	515	1/1	0.96	0.14	86,86,86,86	0
7	CA	E	515	1/1	0.96	0.14	71,71,71,71	0
7	CA	C	507	1/1	0.96	0.24	88,88,88,88	0
7	CA	B	518	1/1	0.96	0.26	98,98,98,98	0
7	CA	C	509	1/1	0.96	0.24	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	MAN	E	507	11/12	0.97	0.11	66,93,115,117	0
7	CA	A	510	1/1	0.97	0.17	127,127,127,127	0
6	MAN	F	504	11/12	0.97	0.13	50,74,97,100	0
7	CA	B	522	1/1	0.97	0.22	66,66,66,66	0
6	MAN	C	505	11/12	0.97	0.15	40,84,99,110	0
7	CA	F	507	1/1	0.97	0.23	77,77,77,77	0
7	CA	B	514	1/1	0.97	0.26	58,58,58,58	0
7	CA	E	516	1/1	0.98	0.22	73,73,73,73	0
7	CA	D	509	1/1	0.98	0.16	69,69,69,69	0
7	CA	E	511	1/1	0.98	0.22	88,88,88,88	0
7	CA	B	521	1/1	0.98	0.24	67,67,67,67	0
7	CA	F	512	1/1	0.98	0.26	68,68,68,68	0
7	CA	E	512	1/1	0.98	0.21	87,87,87,87	0
7	CA	C	514	1/1	0.98	0.19	67,67,67,67	0
7	CA	C	508	1/1	0.98	0.21	61,61,61,61	0
7	CA	C	513	1/1	0.98	0.22	83,83,83,83	0
7	CA	F	514	1/1	0.98	0.23	64,64,64,64	0
7	CA	F	510	1/1	0.98	0.24	58,58,58,58	0
7	CA	C	510	1/1	0.98	0.25	76,76,76,76	0
7	CA	A	508	1/1	0.99	0.21	92,92,92,92	0
7	CA	F	508	1/1	0.99	0.24	62,62,62,62	0
7	CA	F	509	1/1	0.99	0.26	72,72,72,72	0
7	CA	E	514	1/1	0.99	0.20	54,54,54,54	0
7	CA	B	516	1/1	0.99	0.23	64,64,64,64	0
7	CA	C	511	1/1	0.99	0.24	41,41,41,41	0
7	CA	D	513	1/1	0.99	0.26	83,83,83,83	0
7	CA	F	513	1/1	0.99	0.27	57,57,57,57	0
7	CA	D	511	1/1	0.99	0.19	53,53,53,53	0
7	CA	A	515	1/1	0.99	0.20	54,54,54,54	0
7	CA	A	513	1/1	0.99	0.28	75,75,75,75	0
7	CA	E	513	1/1	0.99	0.25	71,71,71,71	0
7	CA	E	517	1/1	0.99	0.18	71,71,71,71	0
7	CA	D	514	1/1	0.99	0.20	77,77,77,77	0
7	CA	B	519	1/1	0.99	0.26	59,59,59,59	0
7	CA	A	516	1/1	0.99	0.20	62,62,62,62	0
7	CA	F	511	1/1	0.99	0.24	54,54,54,54	0
7	CA	C	512	1/1	1.00	0.27	60,60,60,60	0
7	CA	F	506	1/1	1.00	0.25	80,80,80,80	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.