



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 08:17 AM BST

PDB ID : 3FCU  
Title : Structure of headpiece of integrin  $\alpha$ IIBb3 in open conformation  
Authors : Zhu, J.; Luo, B.-H.; Xiao, T.; Zhang, C.; Nishida, N.; Springer, T.A.  
Deposited on : 2008-11-22  
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

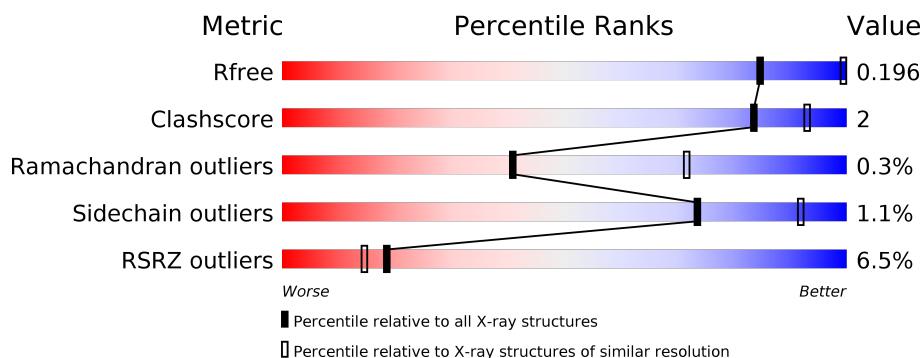
# 1 Overall quality at a glance

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	457	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>8%</div> </div> </div>
1	C	457	<div> <div>2%</div> <div> <div></div> <div>94%</div> <div>• •</div> </div> </div>
1	E	457	<div> <div>%</div> <div> <div></div> <div>93%</div> <div>6% •</div> </div> </div>
2	B	461	<div> <div>9%</div> <div> <div></div> <div>92%</div> <div>7% •</div> </div> </div>
2	D	461	<div> <div>7%</div> <div> <div></div> <div>92%</div> <div>6% •</div> </div> </div>
2	F	461	<div> <div>16%</div> <div> <div></div> <div>91%</div> <div>8% •</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	G	5	 80% 20%
3	I	5	 80% 20%
4	H	4	 75% 25%

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
3	MAN	G	3	X	-	-	X
3	MAN	G	5	-	-	-	X
3	MAN	I	3	X	-	-	X
3	MAN	I	5	-	-	-	X
4	MAN	H	3	X	-	-	X
6	NAG	A	3015	-	-	-	X
6	NAG	C	3015	-	-	-	X
6	NAG	D	3099	-	-	-	X

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 21657 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.

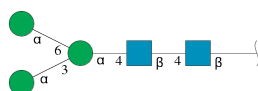
- Molecule 1 is a protein called Integrin, alpha 2b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	0	1	0
			3511	2230	606	667	8			
1	C	451	Total	C	N	O	S	0	2	0
			3477	2207	602	660	8			
1	E	452	Total	C	N	O	S	0	2	0
			3485	2213	603	661	8			

- Molecule 2 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	455	Total	C	N	O	S	0	0	0
			3521	2196	601	692	32			
2	D	455	Total	C	N	O	S	0	0	0
			3521	2196	601	692	32			
2	F	455	Total	C	N	O	S	0	0	0
			3521	2196	601	692	32			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-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
3	G	5	Total	C	N	O	0	0	0
			61	34	2	25			
3	I	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyran

ose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

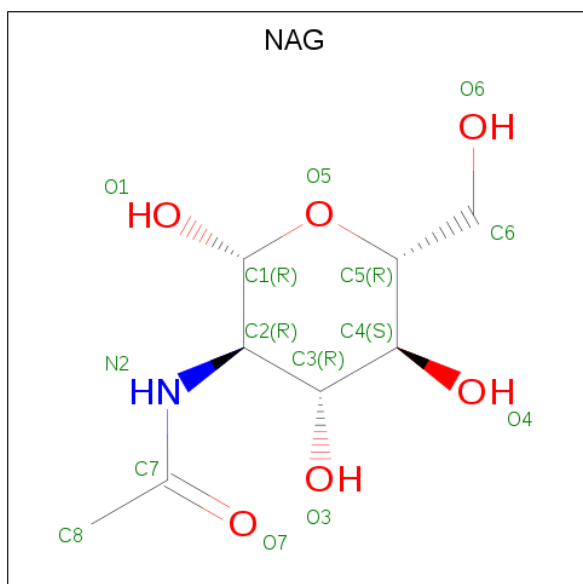


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	2	Total	Ca	0	0
			2	2		
5	E	4	Total	Ca	0	0
			4	4		
5	B	2	Total	Ca	0	0
			2	2		
5	C	4	Total	Ca	0	0
			4	4		
5	A	4	Total	Ca	0	0
			4	4		
5	F	2	Total	Ca	0	0
			2	2		

- Molecule 6 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>).

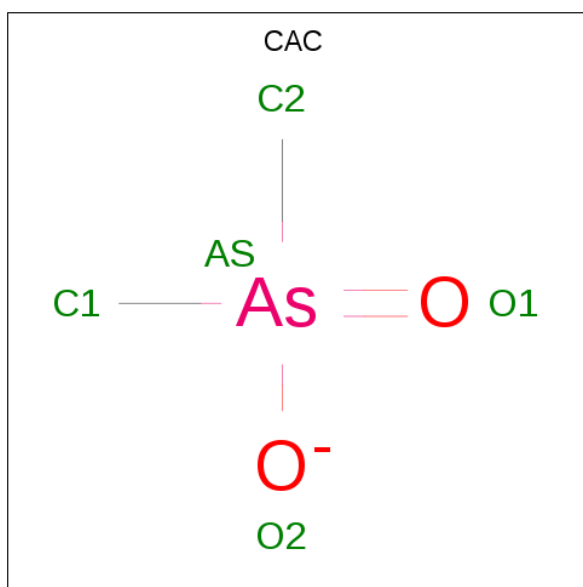


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N O 14 8 1 5	0	0
6	B	1	Total C N O 14 8 1 5	0	0
6	B	1	Total C N O 14 8 1 5	0	0
6	C	1	Total C N O 14 8 1 5	0	0
6	D	1	Total C N O 14 8 1 5	0	0
6	D	1	Total C N O 14 8 1 5	0	0
6	E	1	Total C N O 14 8 1 5	0	0
6	F	1	Total C N O 14 8 1 5	0	0

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	D	1	Total Mg 1 1	0	0
7	E	1	Total Mg 1 1	0	0
7	B	1	Total Mg 1 1	0	0
7	C	1	Total Mg 1 1	0	0
7	A	1	Total Mg 1 1	0	0
7	F	1	Total Mg 1 1	0	0

- Molecule 8 is CACODYLATE ION (three-letter code: CAC) (formula: C<sub>2</sub>H<sub>6</sub>AsO<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	1	Total	As	C	O	0	0
			5	1	2	2		
8	D	1	Total	As	C	O	0	0
			5	1	2	2		
8	F	1	Total	As	C	O	0	0
			5	1	2	2		

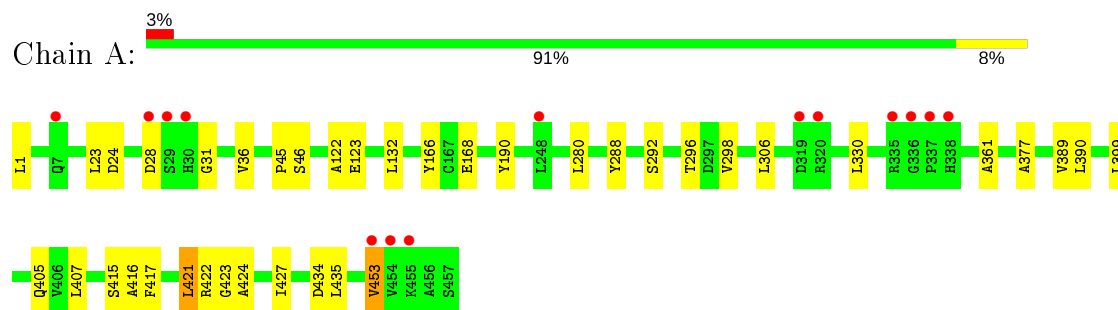
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	191	Total	O	0	0
			191	191		
9	C	56	Total	O	0	0
			56	56		
9	E	51	Total	O	0	0
			51	51		

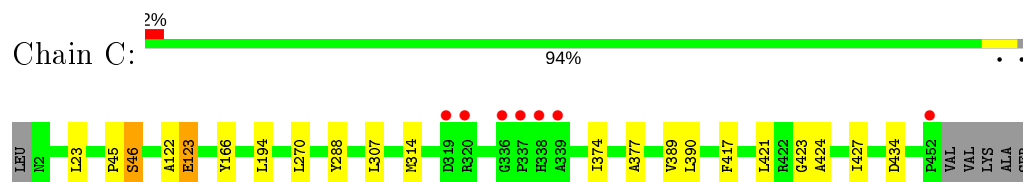
### 3 Residue-property plots [i](#)

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.

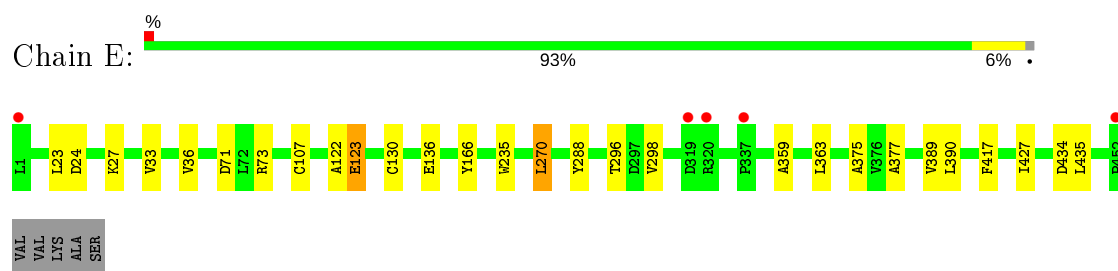
- Molecule 1: Integrin, alpha 2b



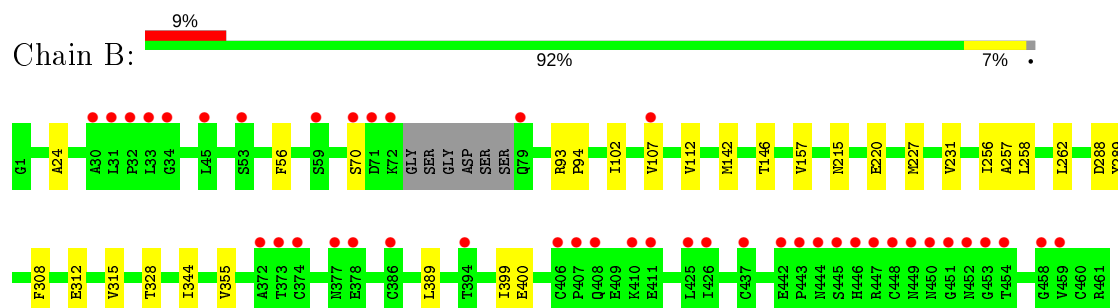
- Molecule 1: Integrin, alpha 2b



- Molecule 1: Integrin, alpha 2b

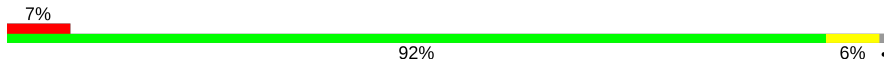


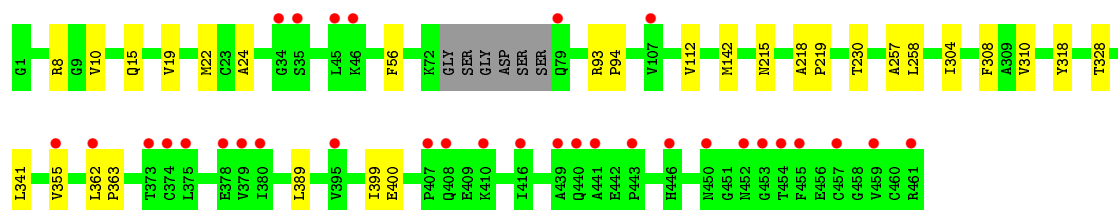
- Molecule 2: Integrin beta-3



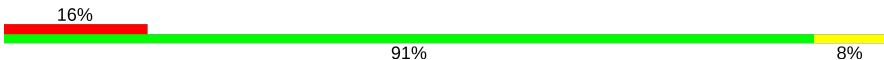


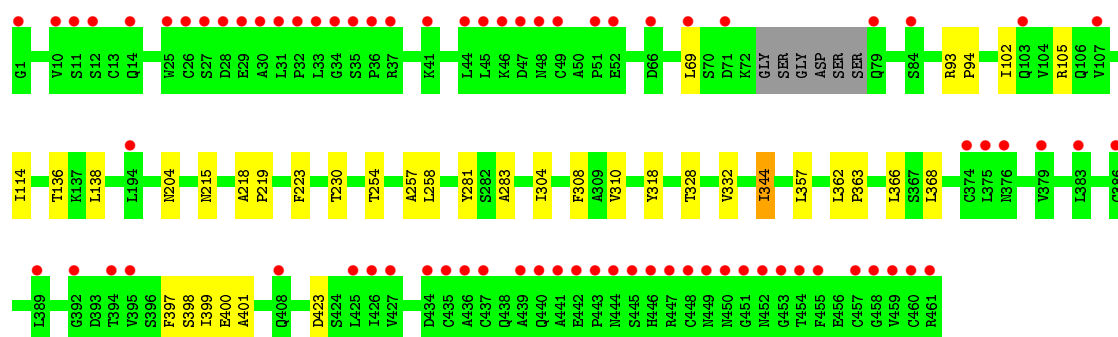
- Molecule 2: Integrin beta-3

Chain D:  7% 92% 6%



- Molecule 2: Integrin beta-3

Chain F:  16% 91% 8%




- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  80% 20%



- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  80% 20%



- Molecule 4: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  75% 25%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	332.09 Å   332.09 Å   88.29 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	44.59 – 2.90 44.58 – 2.89	Depositor EDS
% Data completeness (in resolution range)	95.2 (44.59-2.90) 95.2 (44.58-2.89)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.32 (at 2.90 Å)	Xtriage
Refinement program	REFMAC 5.4.0066	Depositor
R, $R_{free}$	0.174 , 0.197 0.175 , 0.196	Depositor DCC
$R_{free}$ test set	3098 reflections (2.61%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.7	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 72.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	21657	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.65% 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: CAC, MG, 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	0.33	0/3608	0.41	0/4916
1	C	0.32	0/3577	0.41	0/4873
1	E	0.33	0/3585	0.41	0/4884
2	B	0.30	0/3584	0.38	0/4858
2	D	0.30	0/3584	0.39	0/4858
2	F	0.30	0/3584	0.38	0/4858
All	All	0.31	0/21522	0.40	0/29247

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	3511	0	3345	19	0
1	C	3477	0	3303	10	0
1	E	3485	0	3317	15	0
2	B	3521	0	3455	19	0
2	D	3521	0	3455	14	0
2	F	3521	0	3456	23	0
3	G	61	0	52	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	61	0	52	0	0
4	H	50	0	43	0	0
5	A	4	0	0	0	0
5	B	2	0	0	0	0
5	C	4	0	0	0	0
5	D	2	0	0	0	0
5	E	4	0	0	0	0
5	F	2	0	0	0	0
6	A	14	0	13	0	0
6	B	28	0	26	1	0
6	C	14	0	13	0	0
6	D	28	0	26	1	0
6	E	14	0	13	0	0
6	F	14	0	13	1	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
7	E	1	0	0	0	0
7	F	1	0	0	0	0
8	B	5	0	0	0	0
8	D	5	0	0	0	0
8	F	5	0	0	0	0
9	A	191	0	0	0	0
9	C	56	0	0	0	0
9	E	51	0	0	1	0
All	All	21657	0	20582	99	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 2.

All (99) 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)
2:B:112:VAL:HG11	2:B:142:MET:HE3	1.71	0.72
1:A:427:ILE:HG22	1:A:434:ASP:OD1	1.91	0.70
2:F:138:LEU:CD2	2:F:344:ILE:HD12	2.26	0.66
2:D:230:THR:HG23	2:D:304:ILE:HD12	1.79	0.63
2:D:8:ARG:HB2	2:D:10:VAL:HG13	1.80	0.62
2:F:138:LEU:HD23	2:F:344:ILE:HD12	1.84	0.59
1:C:427:ILE:HG22	1:C:434:ASP:OD1	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:427:ILE:HG22	1:E:434:ASP:OD1	2.03	0.59
2:F:400:GLU:HB2	6:F:3371:NAG:H83	1.86	0.58
2:F:223:PHE:CZ	2:F:254:THR:HG21	2.40	0.57
1:E:363:LEU:HD21	1:E:435:LEU:HD13	1.86	0.57
2:D:308:PHE:CE2	2:D:328:THR:HG21	2.40	0.57
1:A:377:ALA:HB2	1:A:421:LEU:HD22	1.86	0.57
2:D:400:GLU:HB2	6:D:3371:NAG:H83	1.86	0.57
2:F:281:TYR:CE1	2:F:283:ALA:HB3	2.41	0.55
1:A:330:LEU:HD22	1:A:399:LEU:HD12	1.89	0.54
2:B:400:GLU:HB2	6:B:3371:NAG:H83	1.90	0.53
2:B:142:MET:HE2	2:B:344:ILE:HG23	1.90	0.52
2:F:218:ALA:HB3	2:F:219:PRO:HD3	1.91	0.52
1:E:122:ALA:O	1:E:123:GLU:HB2	2.08	0.52
1:C:122:ALA:O	1:C:123:GLU:HB2	2.09	0.52
2:B:355:VAL:HG23	2:B:389:LEU:HD22	1.91	0.52
1:A:1:LEU:HD12	1:A:405:GLN:HG3	1.92	0.51
2:D:218:ALA:HB3	2:D:219:PRO:HD3	1.93	0.51
1:C:45:PRO:O	1:C:46:SER:CB	2.57	0.51
2:B:256:ILE:HD12	9:E:494:HOH:O	2.09	0.51
2:B:102:ILE:HG22	2:B:399:ILE:HD11	1.93	0.51
2:F:308:PHE:CE2	2:F:328:THR:HG21	2.45	0.51
1:C:377:ALA:HB2	1:C:421:LEU:HD11	1.93	0.50
1:A:132:LEU:N	1:A:132:LEU:HD12	2.28	0.49
2:F:230:THR:HG23	2:F:304:ILE:HD12	1.93	0.49
2:D:112:VAL:HG21	2:D:142:MET:CE	2.42	0.49
1:E:235:TRP:HZ2	1:E:270:LEU:HD11	1.77	0.49
1:E:296:THR:HG23	1:E:298:VAL:HG13	1.94	0.49
1:C:389:VAL:HG23	1:C:417:PHE:CE2	2.47	0.48
1:C:194:LEU:HD12	1:C:194:LEU:C	2.34	0.48
2:F:138:LEU:HD22	2:F:344:ILE:HD12	1.96	0.48
2:B:24:ALA:HB2	2:B:56:PHE:CD1	2.49	0.47
1:E:24:ASP:O	1:E:36:VAL:HG12	2.13	0.47
1:A:122:ALA:O	1:A:123:GLU:HB2	2.14	0.47
2:F:114:ILE:HD13	2:F:344:ILE:HD11	1.96	0.47
2:D:257:ALA:O	2:D:258:LEU:HB2	2.15	0.47
2:F:399:ILE:HD12	2:F:399:ILE:N	2.30	0.47
2:B:257:ALA:O	2:B:258:LEU:HB2	2.15	0.46
2:B:312:GLU:O	2:B:315:VAL:HG12	2.15	0.46
2:F:332:VAL:O	2:F:332:VAL:HG13	2.14	0.46
1:E:71:ASP:OD1	1:E:73:ARG:HD3	2.16	0.46
2:D:15:GLN:O	2:D:19:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:390:LEU:HD12	1:E:390:LEU:N	2.31	0.46
1:C:390:LEU:HD12	1:C:390:LEU:N	2.31	0.46
2:F:93:ARG:HB2	2:F:94:PRO:HD2	1.98	0.45
1:E:363:LEU:CD2	1:E:435:LEU:HD13	2.46	0.45
1:A:280:LEU:HD11	1:A:306:LEU:HD23	1.99	0.45
2:B:112:VAL:HG11	2:B:142:MET:CE	2.45	0.45
2:B:308:PHE:CE2	2:B:328:THR:HG21	2.52	0.45
1:A:280:LEU:CD1	1:A:306:LEU:HD23	2.46	0.45
2:B:93:ARG:HB2	2:B:94:PRO:HD2	1.99	0.45
2:F:362:LEU:HD12	2:F:363:PRO:HD2	1.97	0.45
2:F:357:LEU:HD11	2:F:397:PHE:CD2	2.51	0.45
2:F:368:LEU:HD23	2:F:401:ALA:HA	1.98	0.44
1:E:363:LEU:HD11	1:E:375:ALA:HB2	1.99	0.44
2:D:355:VAL:HG23	2:D:389:LEU:HD22	2.00	0.44
2:D:93:ARG:HB2	2:D:94:PRO:HD2	1.99	0.44
2:B:70:SER:HB2	2:B:107:VAL:HG11	1.99	0.44
1:E:27:LYS:HG3	1:E:33:VAL:HG22	1.99	0.44
2:F:102:ILE:CG2	2:F:397:PHE:HB2	2.48	0.43
1:A:390:LEU:HD12	1:A:390:LEU:N	2.33	0.43
2:B:142:MET:HE2	2:B:344:ILE:CG2	2.48	0.43
2:B:142:MET:CE	2:B:344:ILE:HG23	2.48	0.43
1:A:24:ASP:O	1:A:36:VAL:HG12	2.18	0.43
2:F:69:LEU:HD13	2:F:105:ARG:HB2	2.00	0.43
2:D:310:VAL:HG11	2:D:318:TYR:CD2	2.53	0.43
1:C:307:LEU:HD11	1:C:374:ILE:HG21	2.01	0.43
1:E:107:CYS:HA	1:E:130:CYS:HA	2.00	0.43
2:F:310:VAL:HG11	2:F:318:TYR:CD2	2.53	0.42
2:B:157:VAL:O	2:B:220:GLU:HB3	2.19	0.42
1:A:415:SER:O	1:A:416:ALA:HB3	2.20	0.42
1:C:314:MET:HB2	1:C:314:MET:HE3	1.94	0.42
2:F:102:ILE:HG22	2:F:399:ILE:HD11	2.01	0.42
2:D:24:ALA:HB2	2:D:56:PHE:CD1	2.55	0.42
2:B:227:MET:O	2:B:231:VAL:HG22	2.19	0.41
1:A:389:VAL:HG23	1:A:417:PHE:CE2	2.55	0.41
1:E:389:VAL:HG23	1:E:417:PHE:CE2	2.55	0.41
1:A:296:THR:HG23	1:A:298:VAL:HG13	2.03	0.41
2:D:399:ILE:HD12	2:D:399:ILE:N	2.34	0.41
1:A:422:ARG:O	1:A:435:LEU:HD12	2.20	0.41
2:D:362:LEU:HD12	2:D:363:PRO:HD2	2.01	0.41
1:A:423:GLY:O	1:A:424:ALA:HB3	2.21	0.41
1:E:359:ALA:HB3	1:E:377:ALA:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:423:GLY:O	1:C:424:ALA:HB3	2.21	0.41
1:E:122:ALA:O	1:E:123:GLU:CB	2.69	0.41
2:F:363:PRO:HG2	2:F:366:LEU:HD12	2.03	0.41
1:A:168:GLU:OE1	2:B:262:LEU:O	2.39	0.40
1:A:377:ALA:HB2	1:A:421:LEU:CD2	2.50	0.40
1:A:361:ALA:HB2	1:A:421:LEU:HB3	2.01	0.40
2:B:288:ASP:OD1	2:B:289:TYR:N	2.54	0.40
2:F:136:THR:HG22	2:F:204:ASN:OD1	2.21	0.40
2:F:257:ALA:O	2:F:258:LEU:HB2	2.21	0.40
1:A:405:GLN:NE2	1:A:407:LEU:HD21	2.36	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	456/457 (100%)	435 (95%)	16 (4%)	5 (1%)	14	42
1	C	451/457 (99%)	436 (97%)	13 (3%)	2 (0%)	34	66
1	E	452/457 (99%)	440 (97%)	11 (2%)	1 (0%)	47	78
2	B	451/461 (98%)	437 (97%)	14 (3%)	0	100	100
2	D	451/461 (98%)	435 (96%)	16 (4%)	0	100	100
2	F	451/461 (98%)	434 (96%)	17 (4%)	0	100	100
All	All	2712/2754 (98%)	2617 (96%)	87 (3%)	8 (0%)	41	71

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	A	28	ASP

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Mol	Chain	Res	Type
1	C	46	SER
1	A	31	GLY
1	E	123	GLU
1	C	123	GLU
1	A	45	PRO
1	A	453	VAL

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	365/364 (100%)	358 (98%)	7 (2%)	57	84
1	C	361/364 (99%)	357 (99%)	4 (1%)	73	92
1	E	362/364 (100%)	357 (99%)	5 (1%)	67	89
2	B	405/409 (99%)	403 (100%)	2 (0%)	88	96
2	D	405/409 (99%)	402 (99%)	3 (1%)	84	95
2	F	405/409 (99%)	401 (99%)	4 (1%)	76	92
All	All	2303/2319 (99%)	2278 (99%)	25 (1%)	73	92

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	166	TYR
1	A	190	TYR
1	A	288	TYR
1	A	292	SER
1	A	421	LEU
1	A	453	VAL
2	B	146	THR
2	B	215	ASN
1	C	23	LEU
1	C	166	TYR
1	C	270	LEU

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Mol	Chain	Res	Type
1	C	288	TYR
2	D	22	MET
2	D	215	ASN
2	D	341	LEU
1	E	23	LEU
1	E	136	GLU
1	E	166	TYR
1	E	270	LEU
1	E	288	TYR
2	F	215	ASN
2	F	344	ILE
2	F	398	SER
2	F	423	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	451	GLN
1	C	2	ASN
2	D	428	GLN
1	E	197	GLN
1	E	395	GLN
2	F	316	ASN
2	F	342	GLN
2	F	428	GLN
2	F	449	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

14 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	$\# Z  > 2$
3	NAG	G	1	3,2	14,14,15	0.53	0	17,19,21	0.94	1 (5%)
3	NAG	G	2	3	14,14,15	0.55	0	17,19,21	0.60	0
3	MAN	G	3	3	11,11,12	0.61	0	15,15,17	0.77	0
3	MAN	G	4	3	11,11,12	0.58	0	15,15,17	0.70	0
3	MAN	G	5	3	11,11,12	0.55	0	15,15,17	0.74	0
4	NAG	H	1	2,4	14,14,15	0.56	0	17,19,21	0.86	1 (5%)
4	NAG	H	2	4	14,14,15	0.45	0	17,19,21	0.90	0
4	MAN	H	3	4	11,11,12	0.59	0	15,15,17	0.61	0
4	MAN	H	4	4	11,11,12	0.58	0	15,15,17	0.65	0
3	NAG	I	1	3,2	14,14,15	0.57	0	17,19,21	0.85	1 (5%)
3	NAG	I	2	3	14,14,15	0.52	0	17,19,21	0.63	0
3	MAN	I	3	3	11,11,12	0.57	0	15,15,17	0.64	0
3	MAN	I	4	3	11,11,12	0.53	0	15,15,17	0.70	0
3	MAN	I	5	3	11,11,12	0.55	0	15,15,17	0.74	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
3	NAG	G	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	MAN	G	3	3	1/1/4/5	2/2/19/22	0/1/1/1
3	MAN	G	4	3	-	1/2/19/22	0/1/1/1
3	MAN	G	5	3	-	2/2/19/22	0/1/1/1
4	NAG	H	1	2,4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	1/6/23/26	0/1/1/1
4	MAN	H	3	4	1/1/4/5	0/2/19/22	0/1/1/1
4	MAN	H	4	4	-	0/2/19/22	0/1/1/1
3	NAG	I	1	3,2	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	MAN	I	3	3	1/1/4/5	0/2/19/22	0/1/1/1
3	MAN	I	4	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	I	5	3	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	NAG	C1-O5-C5	2.94	116.17	112.19
4	H	1	NAG	C1-O5-C5	2.53	115.62	112.19
3	I	1	NAG	C1-O5-C5	2.09	115.03	112.19

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	I	3	MAN	C1
3	G	3	MAN	C1
4	H	3	MAN	C1

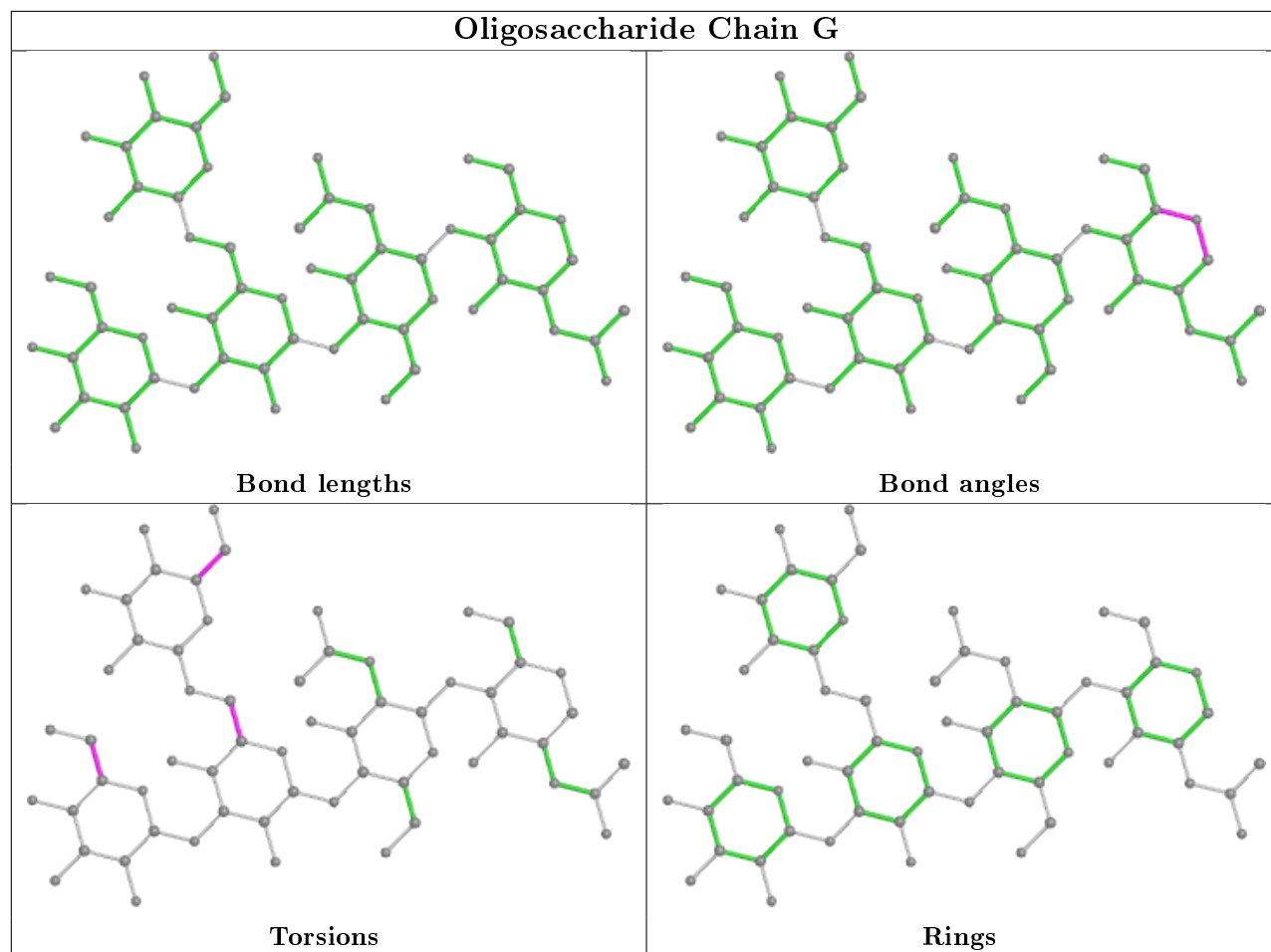
All (10) torsion outliers are listed below:

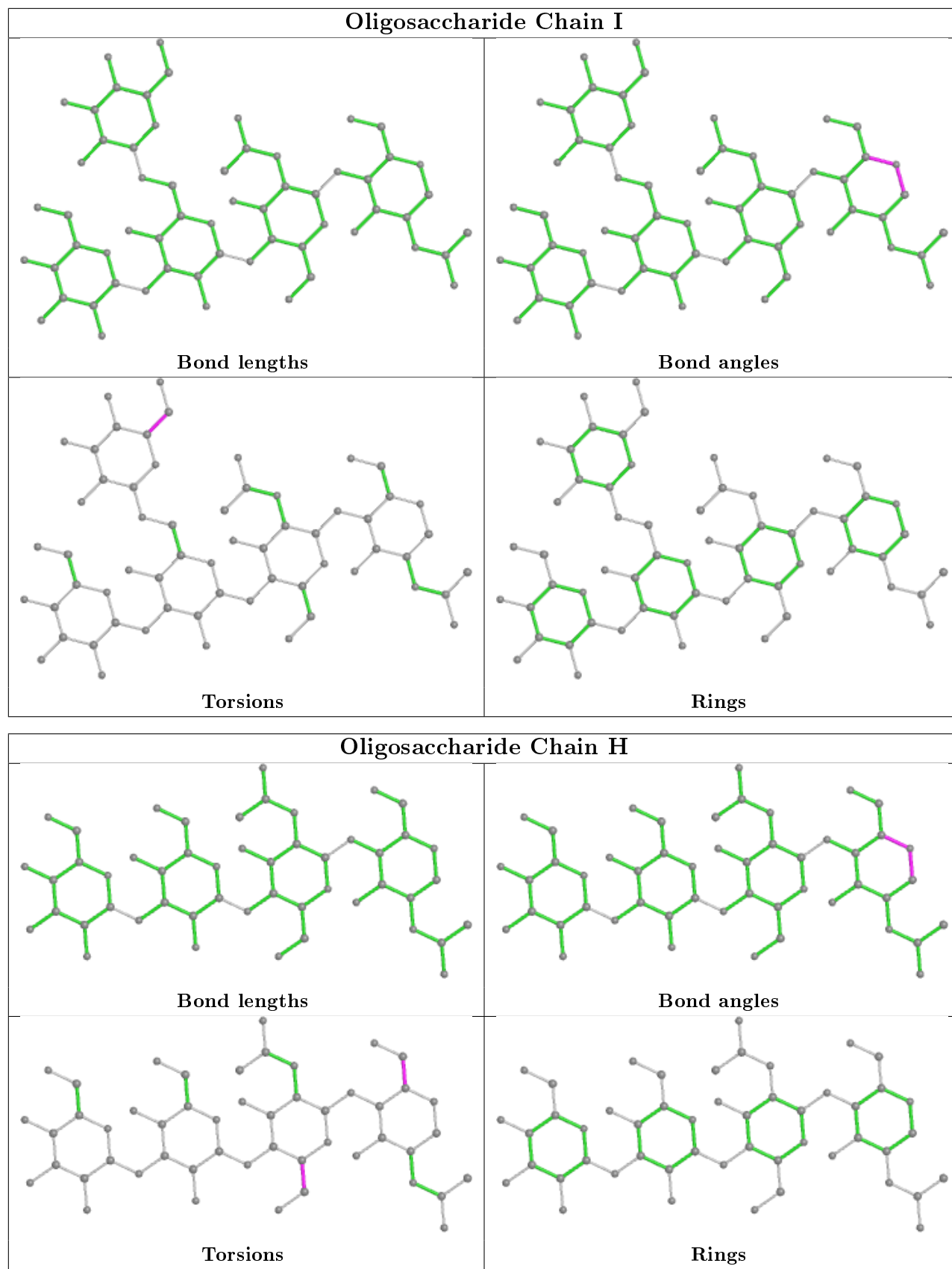
Mol	Chain	Res	Type	Atoms
3	I	5	MAN	O5-C5-C6-O6
3	I	5	MAN	C4-C5-C6-O6
3	G	5	MAN	O5-C5-C6-O6
3	G	5	MAN	C4-C5-C6-O6
4	H	1	NAG	O5-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6
3	G	3	MAN	C4-C5-C6-O6
3	G	3	MAN	O5-C5-C6-O6
3	G	4	MAN	C4-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6

There are no ring outliers.

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 35 ligands modelled in this entry, 24 are monoatomic - leaving 11 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	$\# Z  > 2$
6	NAG	B	3099	2	14,14,15	0.57	0	17,19,21	0.63	0
6	NAG	D	3099	2	14,14,15	0.65	0	17,19,21	0.75	0
6	NAG	F	3371	2	14,14,15	0.52	0	17,19,21	0.68	0
6	NAG	E	3015	1	14,14,15	0.63	0	17,19,21	0.69	0
6	NAG	C	3015	1	14,14,15	0.47	0	17,19,21	0.73	0
8	CAC	B	462	-	0,4,4	0.00	-	0,6,6	0.00	-
8	CAC	D	462	-	0,4,4	0.00	-	0,6,6	0.00	-
6	NAG	A	3015	1	14,14,15	0.60	0	17,19,21	0.65	0
6	NAG	B	3371	2	14,14,15	0.50	0	17,19,21	0.63	0
6	NAG	D	3371	2	14,14,15	0.47	0	17,19,21	0.70	0
8	CAC	F	462	-	0,4,4	0.00	-	0,6,6	0.00	-

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
6	NAG	B	3099	2	-	2/6/23/26	0/1/1/1
6	NAG	D	3099	2	-	2/6/23/26	0/1/1/1
6	NAG	F	3371	2	-	0/6/23/26	0/1/1/1
6	NAG	C	3015	1	-	1/6/23/26	0/1/1/1
6	NAG	E	3015	1	-	2/6/23/26	0/1/1/1
6	NAG	A	3015	1	-	2/6/23/26	0/1/1/1
6	NAG	B	3371	2	-	0/6/23/26	0/1/1/1
6	NAG	D	3371	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	3099	NAG	O5-C5-C6-O6
6	E	3015	NAG	O5-C5-C6-O6
6	A	3015	NAG	O5-C5-C6-O6
6	E	3015	NAG	C4-C5-C6-O6
6	A	3015	NAG	C4-C5-C6-O6
6	B	3099	NAG	C4-C5-C6-O6
6	D	3099	NAG	C4-C5-C6-O6
6	D	3099	NAG	O5-C5-C6-O6
6	C	3015	NAG	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	F	3371	NAG	1	0
6	B	3371	NAG	1	0
6	D	3371	NAG	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, 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	457/457 (100%)	0.33	14 (3%) 49 44	37, 53, 90, 144	0
1	C	451/457 (98%)	0.24	7 (1%) 72 71	37, 52, 85, 133	0
1	E	452/457 (98%)	0.18	5 (1%) 80 80	39, 52, 85, 133	0
2	B	455/461 (98%)	0.61	43 (9%) 8 6	45, 80, 125, 139	1 (0%)
2	D	455/461 (98%)	0.53	32 (7%) 16 12	45, 80, 125, 139	1 (0%)
2	F	455/461 (98%)	0.83	75 (16%) 1 1	45, 80, 124, 139	1 (0%)
All	All	2725/2754 (98%)	0.45	176 (6%) 18 14	37, 60, 122, 144	3 (0%)

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	452	PRO	7.6
2	F	448	CYS	7.1
2	B	453	GLY	6.9
2	B	446	HIS	6.6
2	F	31	LEU	6.4
2	B	452	ASN	6.4
2	F	460	CYS	6.2
2	F	450	ASN	6.1
2	F	441	ALA	6.1
1	C	337	PRO	6.0
2	F	454	THR	5.7
2	F	443	PRO	5.5
2	F	453	GLY	5.3
2	D	446	HIS	5.3
2	B	450	ASN	5.3
2	B	451	GLY	5.2
2	F	48	ASN	5.2
2	F	446	HIS	5.1
2	B	448	CYS	5.1

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Mol	Chain	Res	Type	RSRZ
2	F	461	ARG	4.9
2	F	439	ALA	4.9
2	F	392	GLY	4.9
2	B	30	ALA	4.9
1	A	455	LYS	4.8
2	B	33	LEU	4.7
2	F	375	LEU	4.7
2	B	449	ASN	4.7
1	A	337	PRO	4.6
2	B	410	LYS	4.6
2	B	443	PRO	4.5
1	A	453	VAL	4.4
2	F	49	CYS	4.3
1	A	336	GLY	4.2
2	F	444	ASN	4.2
2	D	375	LEU	4.1
2	F	459	VAL	4.1
2	F	442	GLU	4.1
2	F	28	ASP	4.1
2	F	34	GLY	4.1
2	F	449	ASN	3.9
1	C	338	HIS	3.9
2	F	36	PRO	3.8
2	F	452	ASN	3.8
1	A	28	ASP	3.8
2	F	451	GLY	3.8
2	D	461	ARG	3.8
2	D	450	ASN	3.7
2	F	46	LYS	3.7
1	E	337	PRO	3.7
2	B	454	THR	3.6
1	C	319	ASP	3.6
2	D	443	PRO	3.6
2	D	107	VAL	3.6
2	F	47	ASP	3.6
1	A	7	GLN	3.5
2	F	447	ARG	3.5
2	F	395	VAL	3.5
2	B	31	LEU	3.5
2	B	444	ASN	3.4
1	E	320	ARG	3.4
2	F	435	CYS	3.4

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Mol	Chain	Res	Type	RSRZ
2	F	14	GLN	3.4
2	F	52	GLU	3.3
2	B	445	SER	3.3
2	F	107	VAL	3.3
2	F	29	GLU	3.3
2	F	32	PRO	3.2
2	F	11	SER	3.2
2	F	27	SER	3.2
2	B	408	GLN	3.2
2	F	437	CYS	3.2
1	A	319	ASP	3.2
1	A	454	VAL	3.2
2	D	408	GLN	3.2
2	D	374	CYS	3.2
1	E	319	ASP	3.2
2	F	33	LEU	3.2
2	B	79	GLN	3.1
2	D	378	GLU	3.1
1	A	29	SER	3.1
2	B	458	GLY	3.1
2	D	79	GLN	3.1
2	F	30	ALA	3.1
2	F	45	LEU	3.1
1	C	336	GLY	3.1
2	F	71	ASP	3.1
2	D	439	ALA	3.0
2	F	69	LEU	3.0
1	C	452	PRO	3.0
2	F	37	ARG	3.0
2	D	441	ALA	3.0
1	A	320	ARG	3.0
2	F	10	VAL	3.0
2	B	425	LEU	3.0
1	A	335	ARG	2.9
2	B	447	ARG	2.9
2	B	426	ILE	2.9
2	D	407	PRO	2.9
2	D	440	GLN	2.9
2	B	378	GLU	2.9
2	B	34	GLY	2.9
2	F	458	GLY	2.9
2	F	376	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	374	CYS	2.8
2	D	355	VAL	2.8
2	B	32	PRO	2.8
2	F	79	GLN	2.8
2	B	45	LEU	2.8
1	A	338	HIS	2.8
2	F	445	SER	2.8
1	A	30	HIS	2.8
2	F	436	ALA	2.8
2	D	452	ASN	2.7
2	F	44	LEU	2.7
2	F	12	SER	2.7
2	F	457	CYS	2.6
2	B	71	ASP	2.6
1	A	248	LEU	2.6
2	F	425	LEU	2.6
2	F	84	SER	2.6
2	D	457	CYS	2.6
2	D	455	PHE	2.6
2	F	440	GLN	2.5
2	F	389	LEU	2.5
2	B	107	VAL	2.5
1	E	1	LEU	2.5
2	D	362	LEU	2.5
2	B	373	THR	2.5
2	F	35	SER	2.5
2	D	379	VAL	2.5
2	F	41	LYS	2.5
2	F	394	THR	2.5
2	B	406	CYS	2.5
2	D	410	LYS	2.5
2	B	59	SER	2.4
2	B	459	VAL	2.4
2	F	66	ASP	2.4
2	D	453	GLY	2.4
2	B	372	ALA	2.4
2	B	72	LYS	2.4
2	F	383	LEU	2.4
2	F	455	PHE	2.4
2	D	35	SER	2.4
2	F	379	VAL	2.4
1	C	320	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	442	GLU	2.3
2	B	437	CYS	2.3
2	F	386	CYS	2.3
2	F	1	GLY	2.3
2	B	411	GLU	2.3
2	D	373	THR	2.3
2	F	25	TRP	2.3
2	B	394	THR	2.3
2	B	70	SER	2.2
2	F	434	ASP	2.2
1	C	339	ALA	2.2
2	D	45	LEU	2.2
2	F	51	PRO	2.2
2	F	103	GLN	2.2
2	B	53	SER	2.2
2	F	194	LEU	2.2
2	D	380	ILE	2.2
2	F	26	CYS	2.2
2	D	459	VAL	2.2
2	D	34	GLY	2.1
2	D	454	THR	2.1
2	F	426	ILE	2.1
2	F	427	VAL	2.1
2	D	416	ILE	2.1
2	B	407	PRO	2.1
2	B	386	CYS	2.1
2	D	46	LYS	2.1
2	D	395	VAL	2.1
2	B	377	ASN	2.1
2	B	374	CYS	2.0
2	F	408	GLN	2.0

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

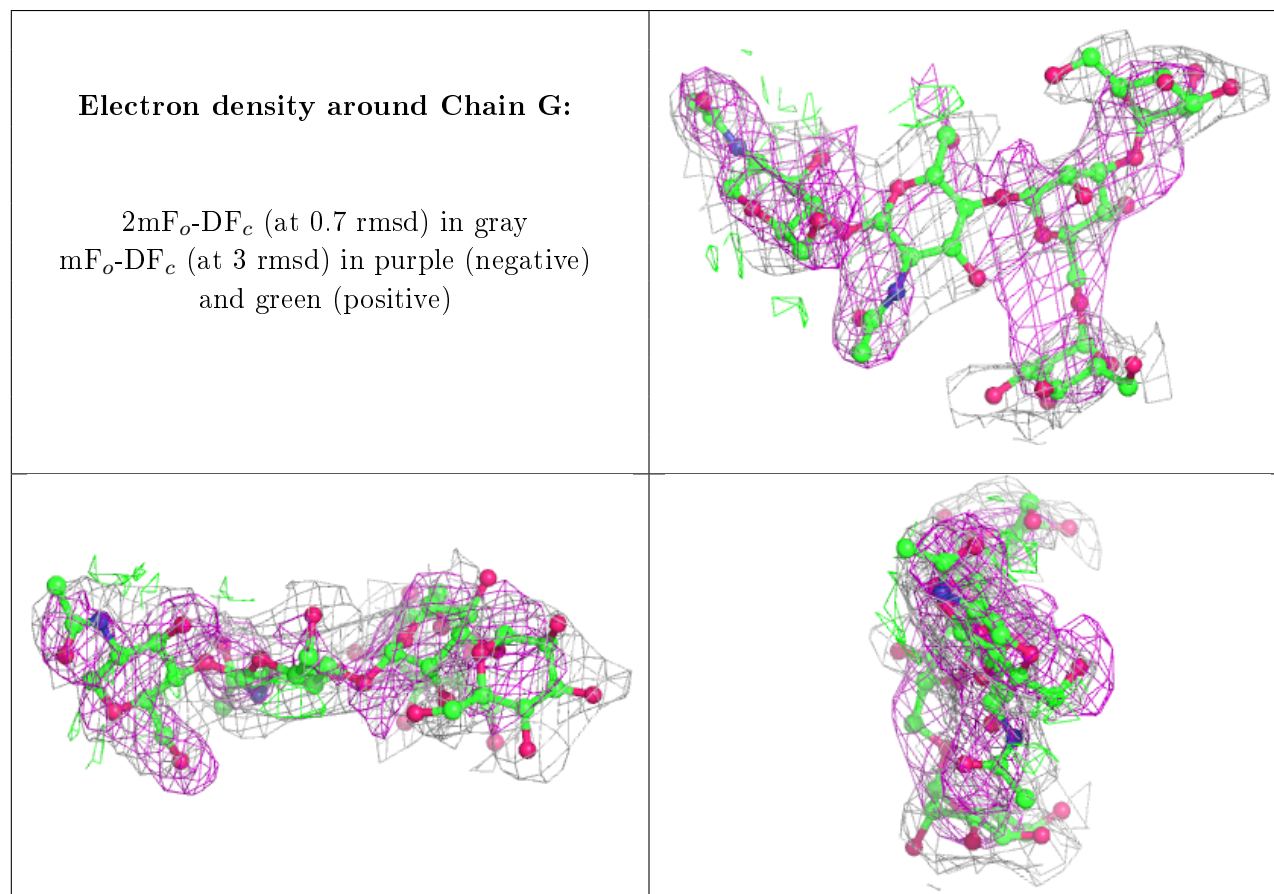
There are no non-standard protein/DNA/RNA residues in this entry.

## 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.

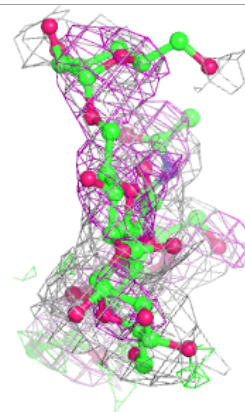
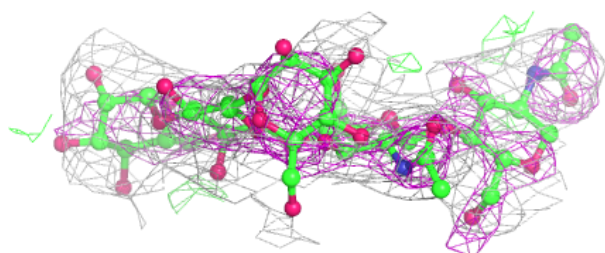
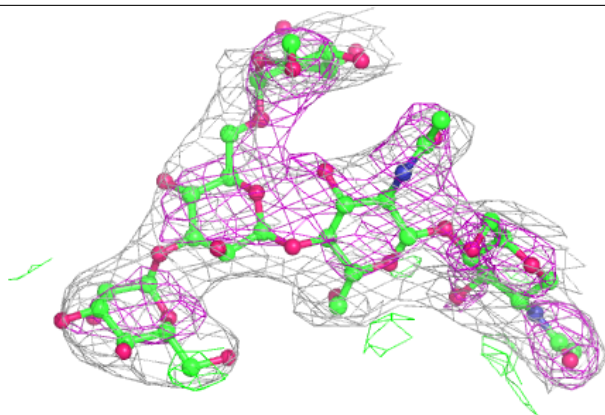
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MAN	G	5	11/12	0.50	0.55	147,148,149,149	0
3	MAN	I	5	11/12	0.59	0.62	144,146,147,147	0
3	MAN	G	3	11/12	0.69	0.50	118,126,137,144	0
4	MAN	H	3	11/12	0.77	0.41	106,114,118,119	0
3	MAN	I	3	11/12	0.78	0.43	118,123,135,142	0
3	MAN	G	4	11/12	0.79	0.39	119,125,127,127	0
4	MAN	H	4	11/12	0.83	0.48	106,113,114,115	0
3	MAN	I	4	11/12	0.88	0.30	111,117,120,121	0
3	NAG	I	2	14/15	0.88	0.34	89,98,102,110	0
3	NAG	G	2	14/15	0.91	0.32	80,86,96,108	0
4	NAG	H	2	14/15	0.92	0.29	89,94,99,103	0
3	NAG	G	1	14/15	0.93	0.23	55,62,65,72	0
4	NAG	H	1	14/15	0.95	0.14	49,59,68,80	0
3	NAG	I	1	14/15	0.96	0.22	52,62,73,78	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.

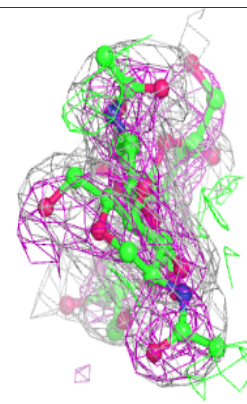
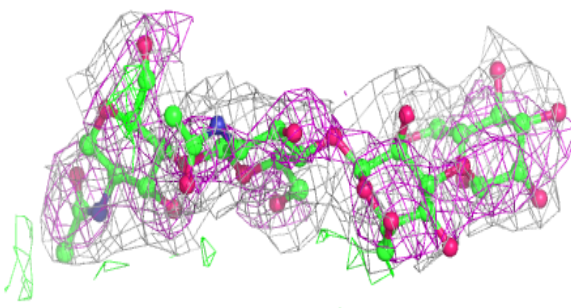
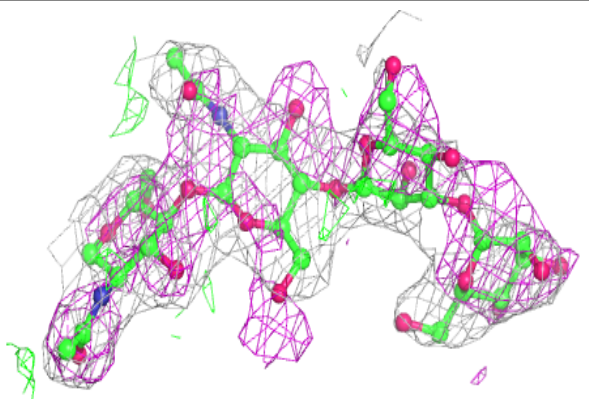


**Electron density around Chain I:**

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

**Electron density around Chain H:**

$2mF_o-DF_c$  (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
6	NAG	A	3015	14/15	0.66	0.51	118,126,129,131	0
6	NAG	D	3099	14/15	0.72	0.44	128,131,135,135	0
6	NAG	B	3099	14/15	0.75	0.40	131,136,139,141	0
6	NAG	D	3371	14/15	0.76	0.24	92,111,115,115	0
6	NAG	E	3015	14/15	0.77	0.33	104,106,108,108	14
6	NAG	C	3015	14/15	0.78	0.45	117,124,126,127	0
6	NAG	F	3371	14/15	0.82	0.38	117,123,126,127	0
6	NAG	B	3371	14/15	0.82	0.39	112,118,120,120	0
7	MG	E	458	1/1	0.91	0.13	68,68,68,68	0
5	CA	B	2002	1/1	0.92	0.09	72,72,72,72	0
5	CA	C	2006	1/1	0.92	0.07	63,63,63,63	0
7	MG	F	2001	1/1	0.93	0.12	37,37,37,37	0
5	CA	C	2007	1/1	0.93	0.04	78,78,78,78	0
5	CA	D	2002	1/1	0.93	0.13	74,74,74,74	0
5	CA	E	2005	1/1	0.94	0.11	42,42,42,42	0
5	CA	A	2005	1/1	0.94	0.08	66,66,66,66	0
5	CA	A	2007	1/1	0.94	0.06	74,74,74,74	0
5	CA	E	2007	1/1	0.94	0.11	65,65,65,65	0
8	CAC	F	462	5/5	0.94	0.23	122,122,124,126	0
8	CAC	D	462	5/5	0.95	0.26	91,96,99,101	0
5	CA	A	2006	1/1	0.95	0.04	61,61,61,61	0
7	MG	D	2001	1/1	0.95	0.12	43,43,43,43	0
8	CAC	B	462	5/5	0.96	0.31	103,106,109,109	0
5	CA	A	2004	1/1	0.97	0.08	56,56,56,56	0
5	CA	D	2003	1/1	0.97	0.22	40,40,40,40	0
7	MG	A	458	1/1	0.97	0.04	61,61,61,61	0
7	MG	C	458	1/1	0.97	0.07	72,72,72,72	0
7	MG	B	2001	1/1	0.97	0.13	39,39,39,39	0
5	CA	E	2004	1/1	0.98	0.10	57,57,57,57	0
5	CA	F	2002	1/1	0.98	0.04	66,66,66,66	0
5	CA	C	2005	1/1	0.98	0.10	43,43,43,43	0
5	CA	C	2004	1/1	0.99	0.07	53,53,53,53	0
5	CA	F	2003	1/1	0.99	0.21	42,42,42,42	0
5	CA	E	2006	1/1	0.99	0.10	62,62,62,62	0
5	CA	B	2003	1/1	1.00	0.21	37,37,37,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.