



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2020 – 03:54 AM BST

PDB ID : 4HSA  
Title : Structure of interleukin 17a in complex with il17ra receptor  
Authors : Liu, S.  
Deposited on : 2012-10-29  
Resolution : 3.15 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

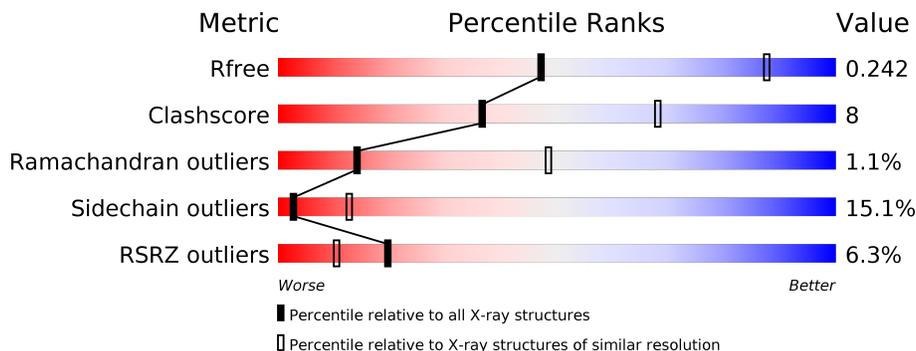
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.15 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	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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	122	
1	B	122	
1	D	122	
1	E	122	
2	C	301	
2	F	301	

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Mol	Chain	Length	Quality of chain
3	G	5	 20% 80%
3	J	5	 40% 60%
4	H	3	 100%
5	I	2	 100%

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	J	4	-	-	-	X
4	FUC	H	3	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8031 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 Interleukin-17A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	104	Total 848	C 532	N 157	O 153	S 6	0	0	0
1	B	109	Total 885	C 547	N 168	O 164	S 6	0	0	0
1	D	100	Total 816	C 511	N 152	O 147	S 6	0	0	0
1	E	109	Total 885	C 547	N 168	O 164	S 6	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	ASP	ASN	engineered mutation	UNP Q16552
A	106	SER	CYS	engineered mutation	UNP Q16552
B	45	ASP	ASN	engineered mutation	UNP Q16552
B	106	SER	CYS	engineered mutation	UNP Q16552
D	45	ASP	ASN	engineered mutation	UNP Q16552
D	106	SER	CYS	engineered mutation	UNP Q16552
E	45	ASP	ASN	engineered mutation	UNP Q16552
E	106	SER	CYS	engineered mutation	UNP Q16552

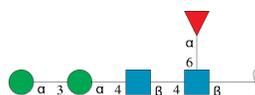
- Molecule 2 is a protein called Interleukin-17 receptor A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	272	Total 2203	C 1388	N 402	O 397	S 16	0	0	0
2	F	271	Total 2193	C 1380	N 401	O 396	S 16	0	0	0

There are 34 discrepancies between the modelled and reference sequences:

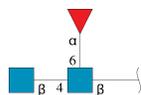
Chain	Residue	Modelled	Actual	Comment	Reference
C	175	ASP	ASN	engineered mutation	UNP Q96F46
C	234	ASP	ASN	engineered mutation	UNP Q96F46
C	287	LEU	-	expression tag	UNP Q96F46
C	288	VAL	-	expression tag	UNP Q96F46
C	289	PRO	-	expression tag	UNP Q96F46
C	290	ARG	-	expression tag	UNP Q96F46
C	291	GLY	-	expression tag	UNP Q96F46
C	292	SER	-	expression tag	UNP Q96F46
C	293	ASP	-	expression tag	UNP Q96F46
C	294	TYR	-	expression tag	UNP Q96F46
C	295	LYS	-	expression tag	UNP Q96F46
C	296	ASP	-	expression tag	UNP Q96F46
C	297	ASP	-	expression tag	UNP Q96F46
C	298	ASP	-	expression tag	UNP Q96F46
C	299	ASP	-	expression tag	UNP Q96F46
C	300	LYS	-	expression tag	UNP Q96F46
C	301	GLY	-	expression tag	UNP Q96F46
F	175	ASP	ASN	engineered mutation	UNP Q96F46
F	234	ASP	ASN	engineered mutation	UNP Q96F46
F	287	LEU	-	expression tag	UNP Q96F46
F	288	VAL	-	expression tag	UNP Q96F46
F	289	PRO	-	expression tag	UNP Q96F46
F	290	ARG	-	expression tag	UNP Q96F46
F	291	GLY	-	expression tag	UNP Q96F46
F	292	SER	-	expression tag	UNP Q96F46
F	293	ASP	-	expression tag	UNP Q96F46
F	294	TYR	-	expression tag	UNP Q96F46
F	295	LYS	-	expression tag	UNP Q96F46
F	296	ASP	-	expression tag	UNP Q96F46
F	297	ASP	-	expression tag	UNP Q96F46
F	298	ASP	-	expression tag	UNP Q96F46
F	299	ASP	-	expression tag	UNP Q96F46
F	300	LYS	-	expression tag	UNP Q96F46
F	301	GLY	-	expression tag	UNP Q96F46

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

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



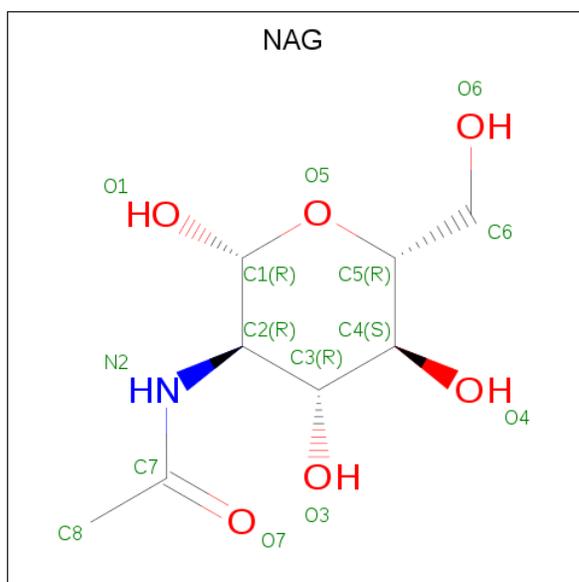
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 5 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- 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
			Total	C	N	O		
6	F	1	14	8	1	5	0	0

- Molecule 7 is water.

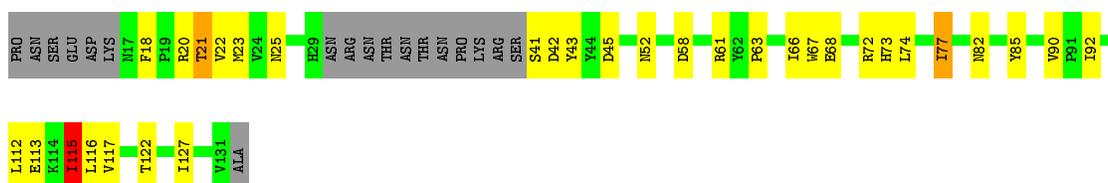
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	2	Total O 2 2	0	0
7	C	1	Total O 1 1	0	0
7	E	2	Total O 2 2	0	0

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

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

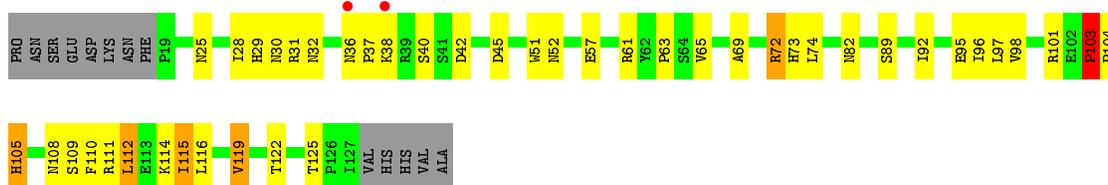
- Molecule 1: Interleukin-17A

Chain A: 



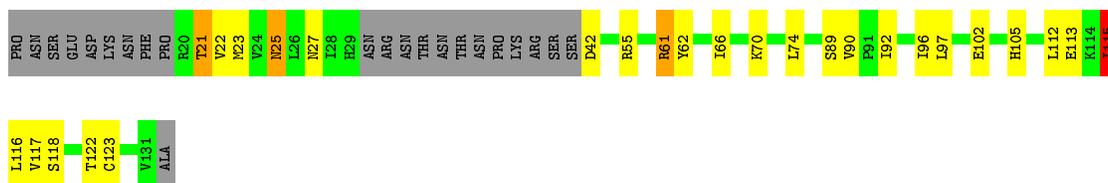
- Molecule 1: Interleukin-17A

Chain B: 



- Molecule 1: Interleukin-17A

Chain D: 



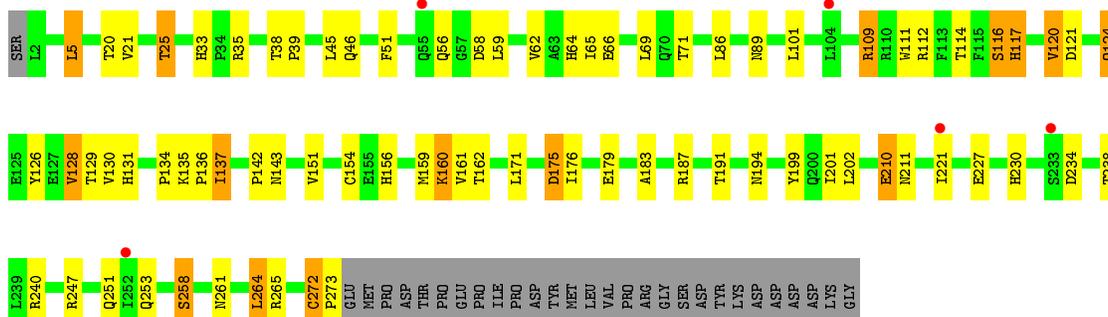
- Molecule 1: Interleukin-17A

Chain E: 

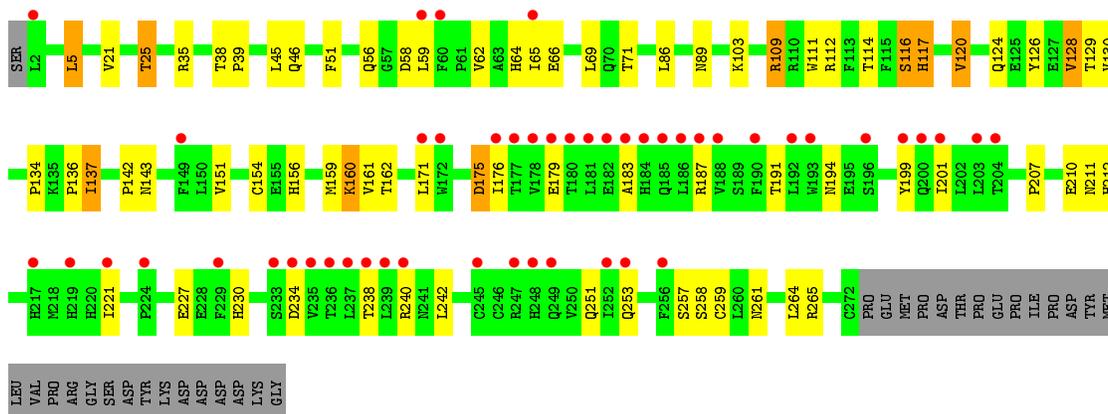




- Molecule 2: Interleukin-17 receptor A



- Molecule 2: Interleukin-17 receptor A



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



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





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

Chain H:  100%



- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.71Å 138.71Å 179.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.14 – 3.15 99.76 – 3.15	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.14-3.15) 100.0 (99.76-3.15)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 3.13Å)	Xtrriage
Refinement program	BUSTER 2.11.2	Depositor
R, $R_{free}$	0.187 , 0.232 0.205 , 0.242	Depositor DCC
$R_{free}$ test set	1776 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	86.7	Xtrriage
Anisotropy	0.459	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 83.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	8031	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	106.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, 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	0.63	1/872 (0.1%)	0.83	0/1190
1	B	0.58	0/908	0.91	3/1238 (0.2%)
1	D	0.54	1/838 (0.1%)	0.77	0/1143
1	E	0.55	0/908	0.91	3/1238 (0.2%)
2	C	0.52	0/2271	0.77	1/3098 (0.0%)
2	F	0.47	0/2260	0.76	1/3082 (0.0%)
All	All	0.53	2/8057 (0.0%)	0.81	8/10989 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	115	ILE	CG1-CD1	5.95	1.91	1.50
1	D	115	ILE	CG1-CD1	5.02	1.85	1.50

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	103	PRO	CA-C-N	6.01	133.94	117.10
1	B	104	PRO	C-N-CA	5.98	136.65	121.70
1	B	119	VAL	N-CA-CB	-5.87	98.59	111.50
2	F	25	THR	N-CA-C	-5.68	95.65	111.00
2	C	25	THR	N-CA-C	-5.41	96.38	111.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	848	0	814	27	0
1	B	885	0	858	19	0
1	D	816	0	787	22	0
1	E	885	0	858	20	0
2	C	2203	0	2081	34	0
2	F	2193	0	2066	25	0
3	G	60	0	52	0	0
3	J	60	0	52	0	0
4	H	38	0	34	2	0
5	I	24	0	22	0	0
6	F	14	0	13	0	0
7	B	2	0	0	0	0
7	C	1	0	0	0	0
7	E	2	0	0	0	0
All	All	8031	0	7637	124	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 8.

The worst 5 of 124 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:D:66:ILE:CG1	1:D:66:ILE:CD1	1.75	1.65
1:A:66:ILE:CD1	1:A:66:ILE:CG1	1.76	1.58
1:E:96:ILE:CG1	1:E:96:ILE:CD1	1.84	1.52
1:D:115:ILE:CG1	1:D:115:ILE:CD1	1.85	1.51
1:A:115:ILE:CD1	1:A:115:ILE:CG1	1.91	1.47

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	100/122 (82%)	93 (93%)	7 (7%)	0	100	100
1	B	107/122 (88%)	98 (92%)	6 (6%)	3 (3%)	5	26
1	D	96/122 (79%)	94 (98%)	2 (2%)	0	100	100
1	E	107/122 (88%)	99 (92%)	5 (5%)	3 (3%)	5	26
2	C	270/301 (90%)	251 (93%)	17 (6%)	2 (1%)	22	59
2	F	269/301 (89%)	253 (94%)	14 (5%)	2 (1%)	22	59
All	All	949/1090 (87%)	888 (94%)	51 (5%)	10 (1%)	14	48

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	40	SER
1	B	103	PRO
1	B	105	HIS
1	E	40	SER
1	E	103	PRO

### 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	99/116 (85%)	89 (90%)	10 (10%)	7	27
1	B	104/116 (90%)	85 (82%)	19 (18%)	1	8
1	D	95/116 (82%)	82 (86%)	13 (14%)	3	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	104/116 (90%)	90 (86%)	14 (14%)	4	17
2	C	251/284 (88%)	212 (84%)	39 (16%)	2	11
2	F	249/284 (88%)	208 (84%)	41 (16%)	2	10
All	All	902/1032 (87%)	766 (85%)	136 (15%)	3	13

5 of 136 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	227	GLU
1	D	102	GLU
2	F	210	GLU
2	C	238	THR
1	D	25	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	131	HIS
1	D	27	ASN
2	F	131	HIS
2	C	156	HIS
2	C	219	HIS

### 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](#)

15 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	0.30	0	17,19,21	0.72	1 (5%)
3	NAG	G	2	3	14,14,15	0.31	0	17,19,21	0.54	0
3	MAN	G	3	3	11,11,12	0.33	0	15,15,17	0.85	1 (6%)
3	MAN	G	4	3	11,11,12	0.39	0	15,15,17	0.93	2 (13%)
3	FUC	G	5	3	10,10,11	0.40	0	14,14,16	0.73	1 (7%)
4	NAG	H	1	2,4	14,14,15	0.42	0	17,19,21	1.50	2 (11%)
4	NAG	H	2	4	14,14,15	0.33	0	17,19,21	0.93	2 (11%)
4	FUC	H	3	4	10,10,11	0.58	0	14,14,16	1.32	2 (14%)
5	NAG	I	1	2,5	14,14,15	0.33	0	17,19,21	0.79	1 (5%)
5	FUC	I	2	5	10,10,11	0.47	0	14,14,16	1.11	1 (7%)
3	NAG	J	1	3,2	14,14,15	0.31	0	17,19,21	0.82	1 (5%)
3	NAG	J	2	3	14,14,15	0.30	0	17,19,21	0.61	0
3	MAN	J	3	3	11,11,12	0.33	0	15,15,17	0.81	1 (6%)
3	MAN	J	4	3	11,11,12	0.43	0	15,15,17	0.93	2 (13%)
3	FUC	J	5	3	10,10,11	0.41	0	14,14,16	0.66	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	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	MAN	G	3	3	-	1/2/19/22	1/1/1/1
3	MAN	G	4	3	-	2/2/19/22	0/1/1/1
3	FUC	G	5	3	-	-	0/1/1/1
4	NAG	H	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	H	2	4	-	0/6/23/26	0/1/1/1
4	FUC	H	3	4	-	-	0/1/1/1
5	NAG	I	1	2,5	-	2/6/23/26	0/1/1/1
5	FUC	I	2	5	-	-	0/1/1/1
3	NAG	J	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	J	2	3	-	2/6/23/26	0/1/1/1
3	MAN	J	3	3	-	1/2/19/22	1/1/1/1

*Continued on next page...*

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	J	4	3	-	2/2/19/22	0/1/1/1
3	FUC	J	5	3	-	-	0/1/1/1

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1	NAG	C1-O5-C5	4.75	118.63	112.19
4	H	3	FUC	C1-O5-C5	4.13	122.14	112.78
4	H	1	NAG	C1-C2-N2	3.35	116.21	110.49
5	I	2	FUC	C1-O5-C5	3.11	119.83	112.78
3	G	3	MAN	C1-O5-C5	2.88	116.10	112.19

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

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

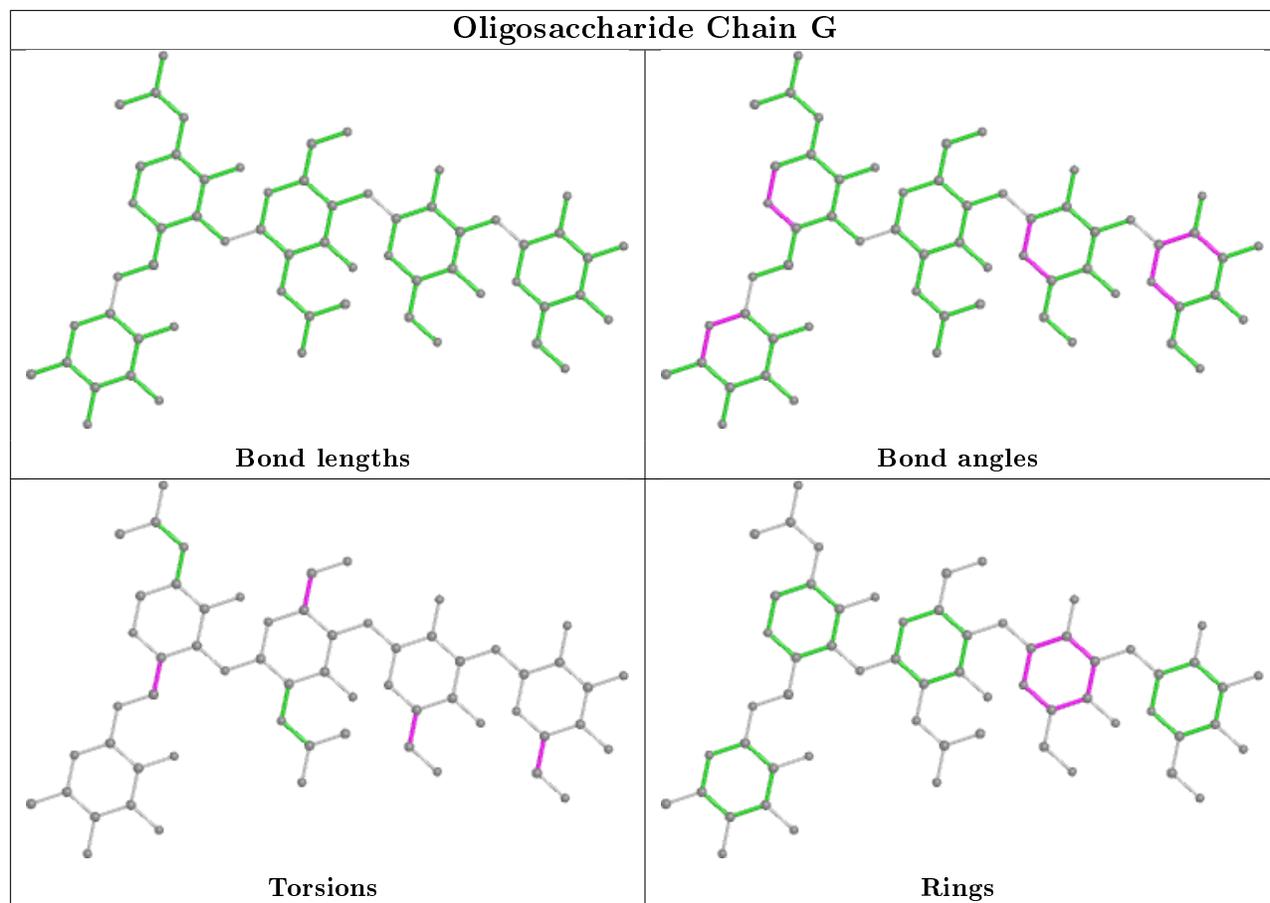
All (2) ring outliers are listed below:

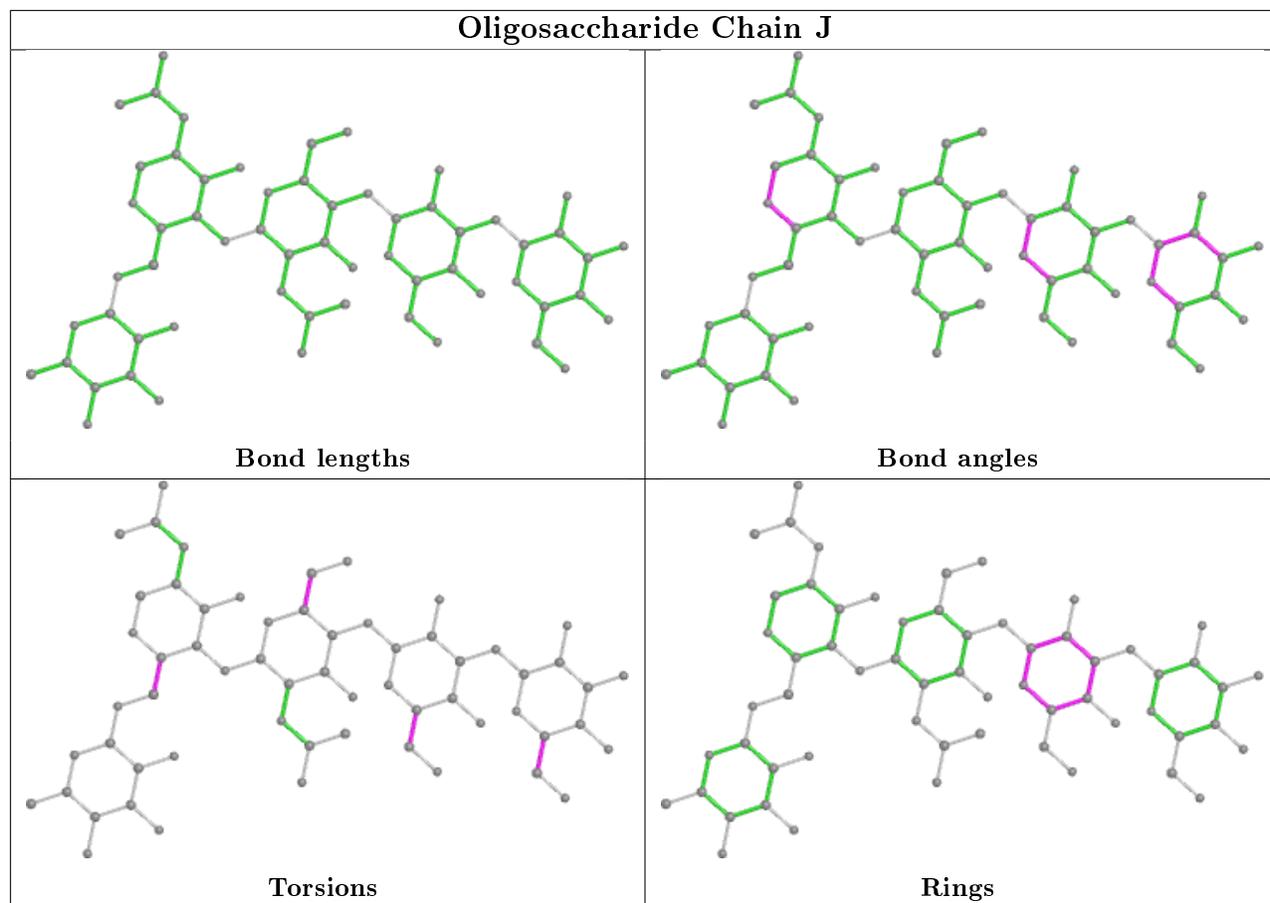
Mol	Chain	Res	Type	Atoms
3	J	3	MAN	C1-C2-C3-C4-C5-O5
3	G	3	MAN	C1-C2-C3-C4-C5-O5

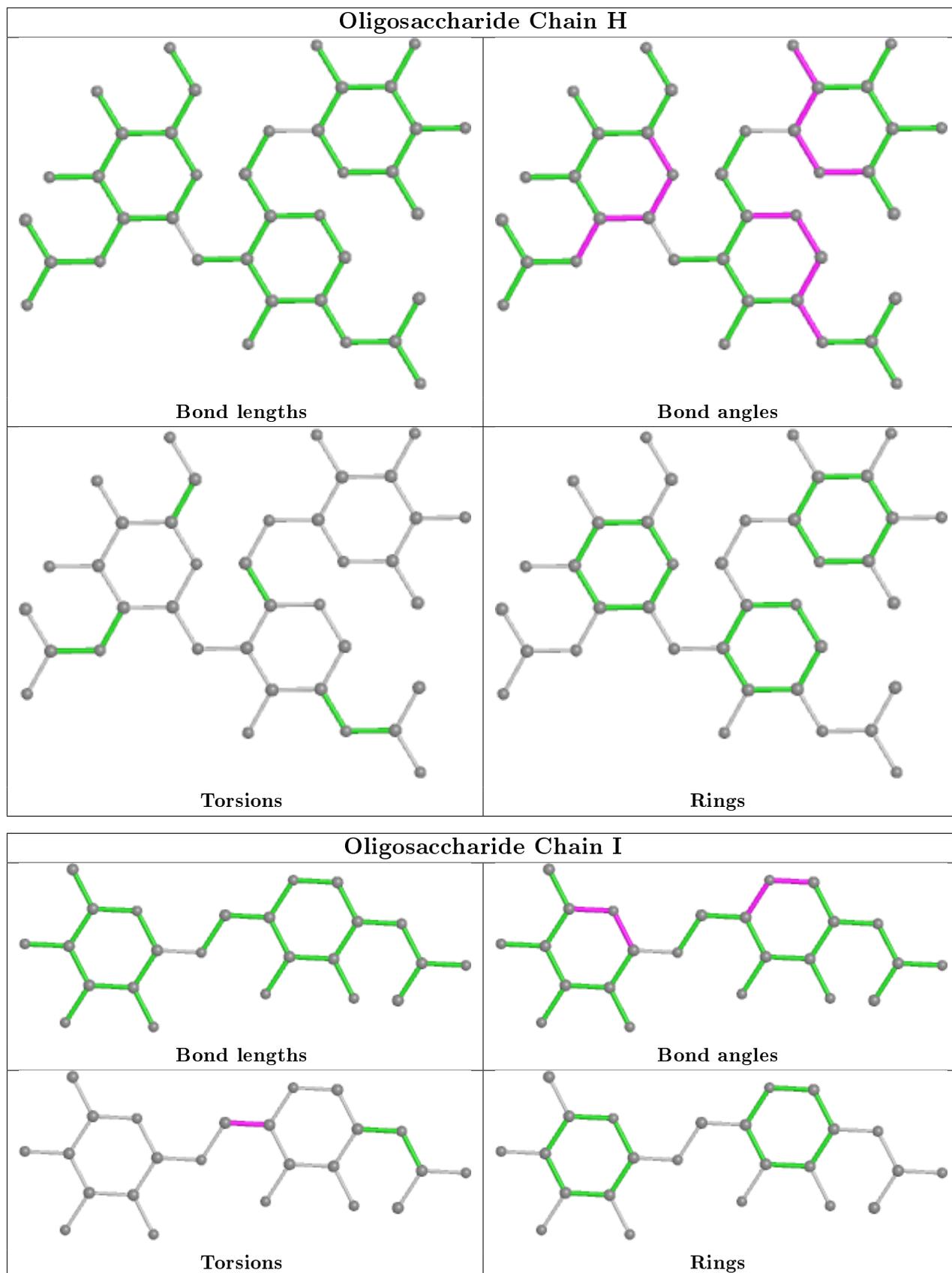
3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	2	NAG	1	0
4	H	1	NAG	2	0
4	H	3	FUC	1	0

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







## 5.6 Ligand geometry [i](#)

1 ligand is 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
6	NAG	F	606	2	14,14,15	0.37	0	17,19,21	1.35	2 (11%)

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	F	606	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
6	F	606	NAG	C1-O5-C5	4.03	117.65	112.19
6	F	606	NAG	C1-C2-N2	3.36	116.22	110.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	104/122 (85%)	0.40	0 <b>100</b> <b>100</b>	59, 87, 121, 158	0
1	B	109/122 (89%)	0.51	2 (1%) 68 55	59, 79, 146, 172	0
1	D	100/122 (81%)	0.21	0 <b>100</b> <b>100</b>	70, 100, 140, 174	0
1	E	109/122 (89%)	0.59	5 (4%) 32 18	58, 83, 139, 183	0
2	C	272/301 (90%)	0.54	5 (1%) 68 55	55, 91, 137, 173	0
2	F	271/301 (90%)	1.06	49 (18%) 1 1	59, 114, 219, 230	0
All	All	965/1090 (88%)	0.64	61 (6%) 20 10	55, 93, 188, 230	0

The worst 5 of 61 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	38	LYS	6.4
2	F	190	PHE	5.2
2	F	171	LEU	5.1
2	F	249	GLN	5.1
2	F	193	TRP	5.1

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

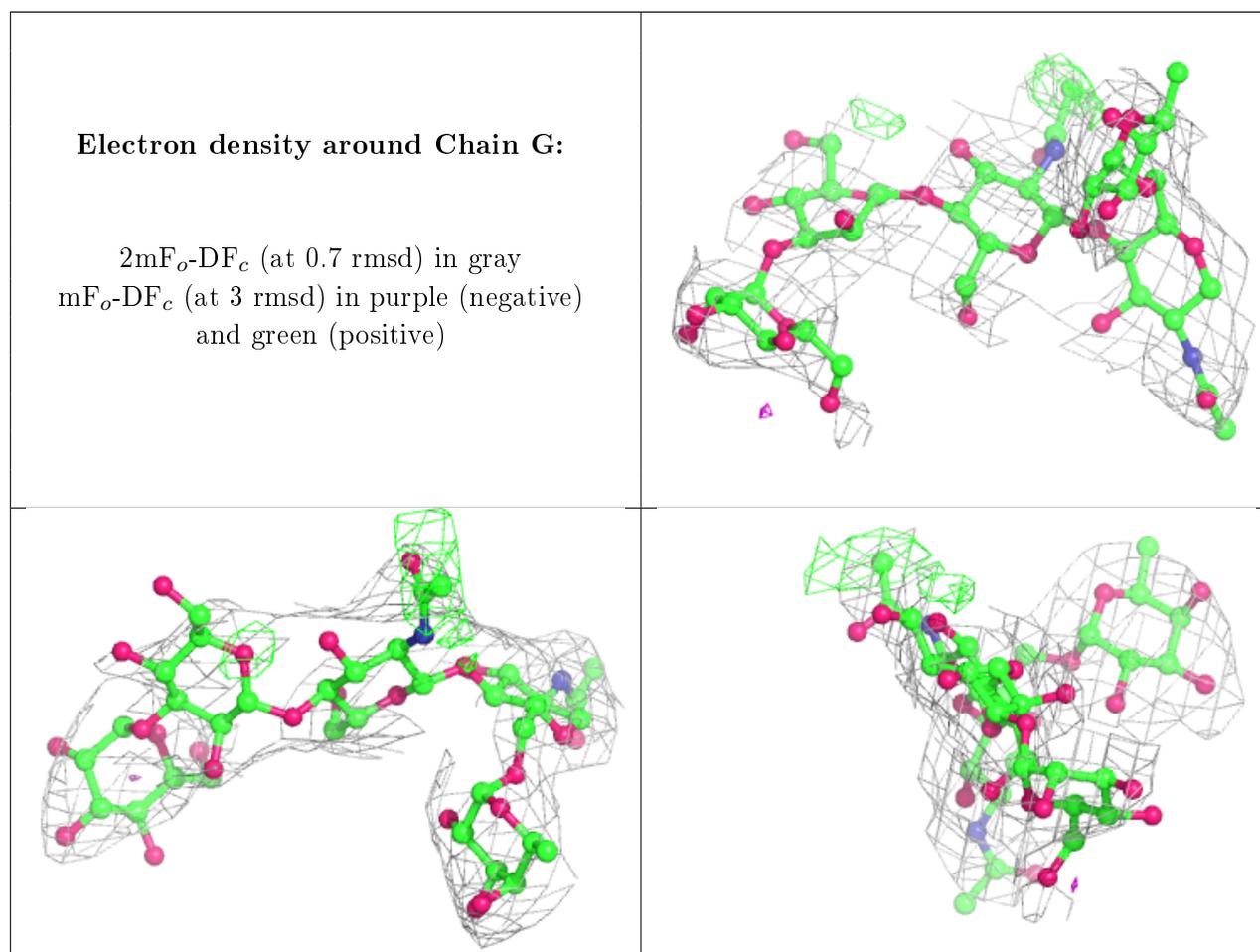
There are no non-standard protein/DNA/RNA residues in this entry.

### 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.

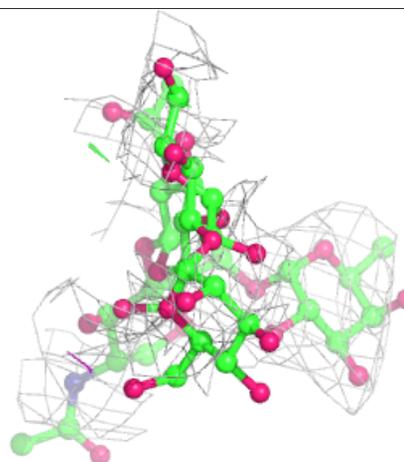
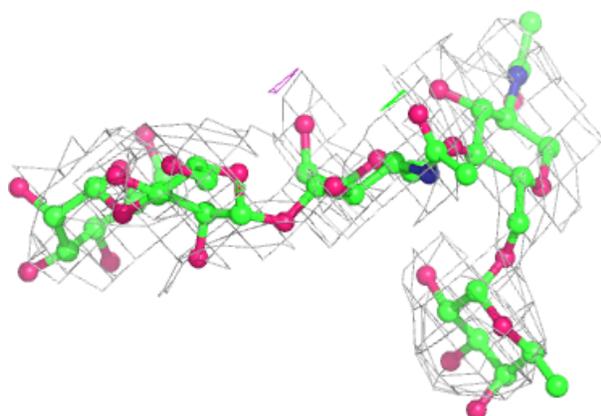
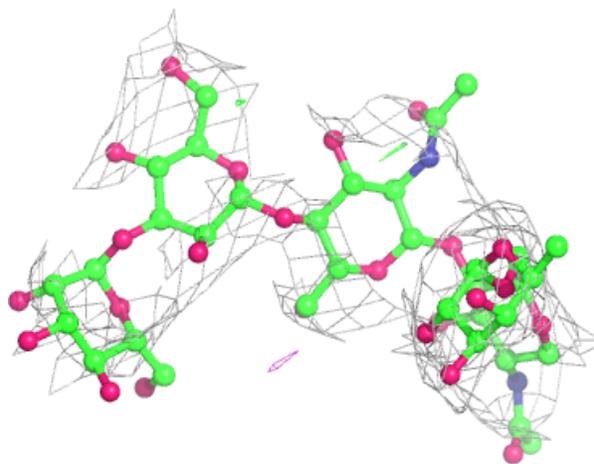
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	NAG	H	1	14/15	0.44	0.24	132,148,162,168	0
3	MAN	G	4	11/12	0.73	0.35	197,199,200,200	0
3	NAG	J	2	14/15	0.76	0.35	196,207,211,214	0
5	FUC	I	2	10/11	0.76	0.34	166,174,177,177	0
4	NAG	H	2	14/15	0.77	0.27	151,161,166,167	0
3	MAN	J	4	11/12	0.77	0.44	225,227,230,231	0
4	FUC	H	3	10/11	0.78	0.90	164,170,174,174	0
3	FUC	J	5	10/11	0.78	0.39	203,204,205,206	0
3	MAN	J	3	11/12	0.79	0.17	213,216,220,223	0
3	MAN	G	3	11/12	0.80	0.24	194,198,199,201	0
3	NAG	J	1	14/15	0.81	0.26	198,202,203,203	0
5	NAG	I	1	14/15	0.83	0.40	149,154,166,166	0
3	NAG	G	1	14/15	0.91	0.16	150,159,164,168	0
3	NAG	G	2	14/15	0.91	0.19	170,180,188,192	0
3	FUC	G	5	10/11	0.92	0.26	162,165,168,171	0

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



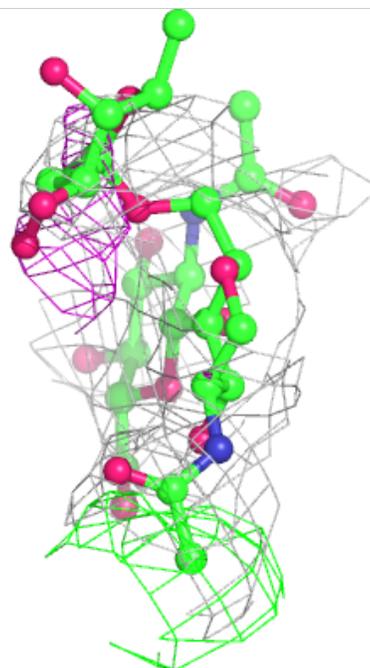
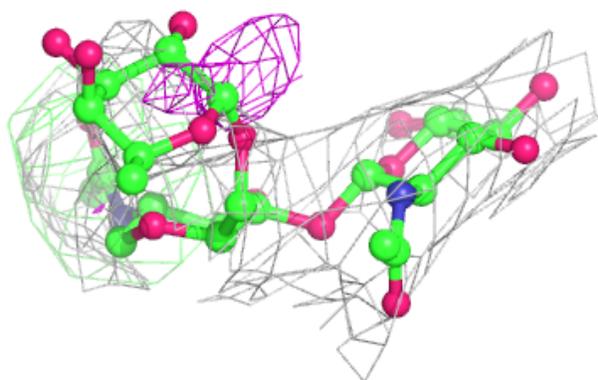
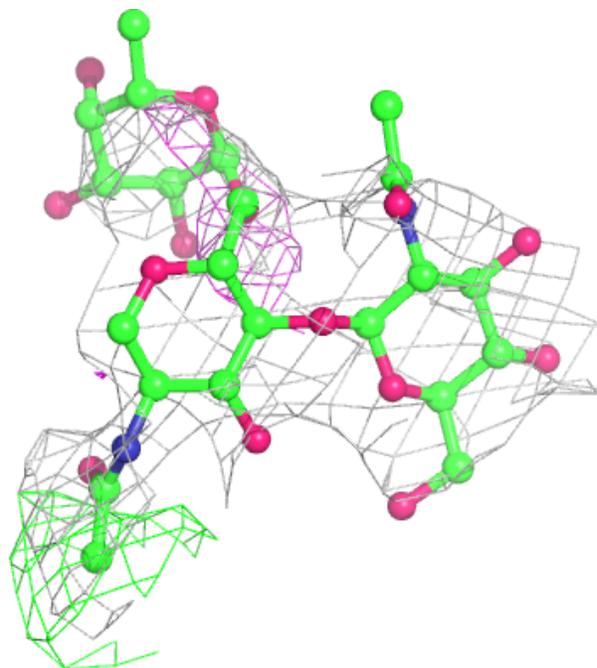
**Electron density around Chain J:**

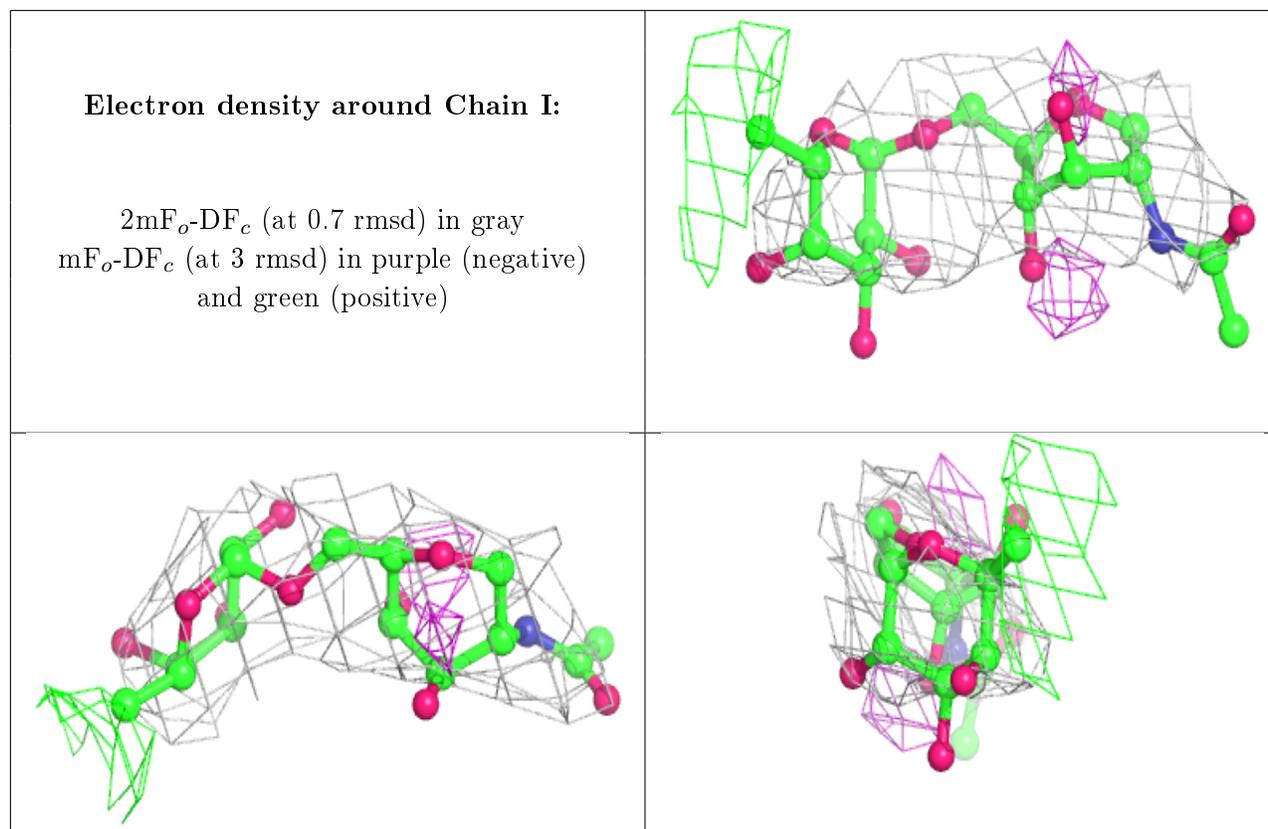
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain 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 [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
6	NAG	F	606	14/15	0.66	0.19	155,163,169,173	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.