



# wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 12, 2023 – 05:55 PM EDT

PDB ID : 4OGA  
Title : Insulin in complex with Site 1 of the human insulin receptor  
Authors : Lawrence, M.C.; Menting, J.G.  
Deposited on : 2014-01-15  
Resolution : 3.50 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)

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<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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)  
Xtrriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

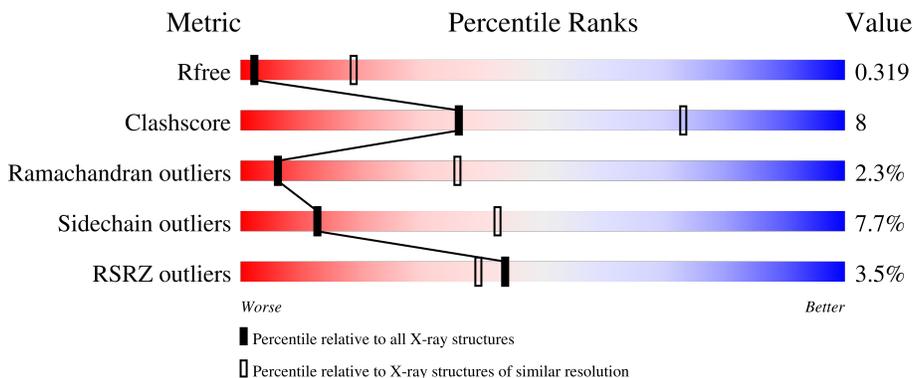
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	21	 5% 38% 48% 14%
2	B	30	 3% 43% 23% 30%
3	C	118	 8% 77% 21% .
4	D	114	 5% 72% 25% .
5	E	317	 % 79% 11% . 9%

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Mol	Chain	Length	Quality of chain
6	F	16	
7	G	4	
8	H	4	

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
8	BMA	H	3	-	-	-	X

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 4693 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 Insulin A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	21	162	99	25	34	4	0	0	0

- Molecule 2 is a protein called Insulin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	21	165	108	26	29	2	0	0	0

- Molecule 3 is a protein called monoclonal antibody fab 83-7 fragment - heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	118	904	568	153	178	5	0	0	0

- Molecule 4 is a protein called monoclonal antibody fab 83-7 fragment - light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	114	891	565	149	173	4	0	0	0

- Molecule 5 is a protein called Insulin receptor domains L1-CR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	288	2301	1449	397	423	32	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	144	HIS	TYR	variant	UNP P06213
E	311	SER	-	expression tag	UNP P06213

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Chain	Residue	Modelled	Actual	Comment	Reference
E	312	SER	-	expression tag	UNP P06213
E	313	SER	-	expression tag	UNP P06213
E	314	LEU	-	expression tag	UNP P06213
E	315	VAL	-	expression tag	UNP P06213
E	316	PRO	-	expression tag	UNP P06213
E	317	ARG	-	expression tag	UNP P06213

- Molecule 6 is a protein called Insulin receptor alpha-CT peptide.

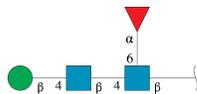
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	F	15	129	86	21	22	0	0	0

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



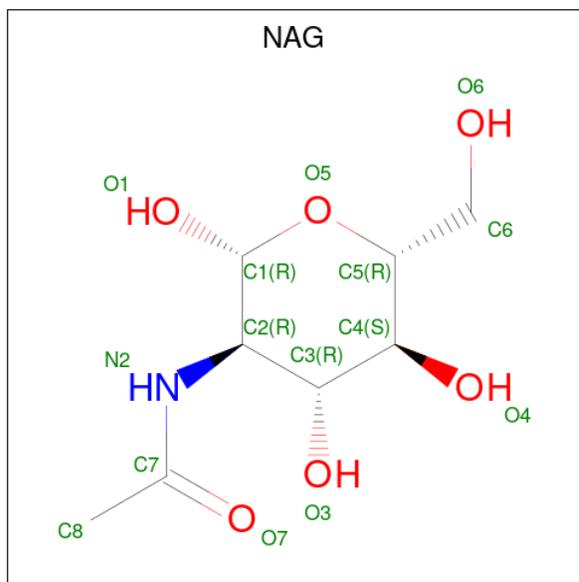
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
7	G	4	50	28	2	20	0	0	0

- Molecule 8 is an oligosaccharide called 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
			Total	C	N	O			
8	H	4	49	28	2	19	0	0	0

- Molecule 9 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
9	E	1	Total	C	N	O	0	0
			14	8	1	5		
9	E	1	Total	C	N	O	0	0
			14	8	1	5		
9	E	1	Total	C	N	O	0	0
			14	8	1	5		

### 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.

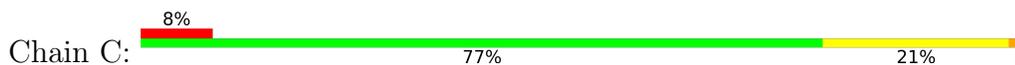
- Molecule 1: Insulin A chain



- Molecule 2: Insulin B chain



- Molecule 3: monoclonal antibody fab 83-7 fragment - heavy chain

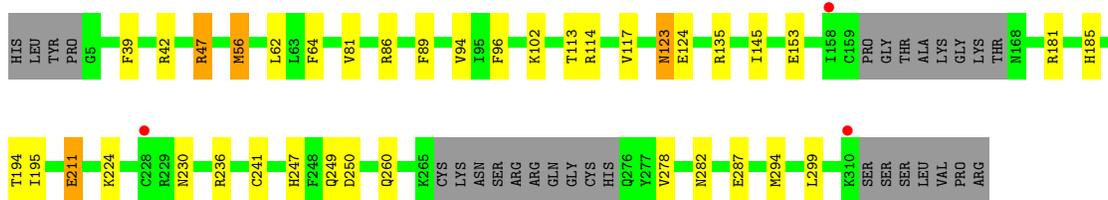


- Molecule 4: monoclonal antibody fab 83-7 fragment - light chain

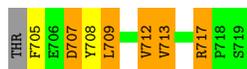


- Molecule 5: Insulin receptor domains L1-CR





- Molecule 6: Insulin receptor alpha-CT peptide



- Molecule 7: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 8: 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



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	169.04Å 169.04Å 169.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 3.50 19.65 – 3.50	Depositor EDS
% Data completeness (in resolution range)	93.3 (19.78-3.50) 93.6 (19.65-3.50)	Depositor EDS
$R_{merge}$	0.21	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.82 (at 3.52Å)	Xtrriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.264 , 0.284 0.283 , 0.319	Depositor DCC
$R_{free}$ test set	942 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	212.0	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 195.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.056 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4693	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	238.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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: MAN, NAG, BMA, FUC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.68	0/163	1.05	0/220
2	B	0.61	0/169	0.86	0/227
3	C	0.49	0/926	0.78	0/1258
4	D	0.52	0/910	0.73	1/1228 (0.1%)
5	E	0.51	0/2355	0.80	0/3192
6	F	0.66	0/134	0.91	0/183
All	All	0.52	0/4657	0.80	1/6308 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	82	SER	C-N-CA	5.24	134.80	121.70

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	162	0	149	7	0
2	B	165	0	152	10	0
3	C	904	0	879	28	0
4	D	891	0	882	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	2301	0	2201	20	0
6	F	129	0	119	3	0
7	G	50	0	43	0	0
8	H	49	0	43	0	0
9	E	42	0	39	0	0
All	All	4693	0	4507	74	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 74 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)
3:C:99:PRO:HG2	3:C:103:LYS:HB2	1.45	0.94
2:B:24:PHE:HE2	5:E:39:PHE:HE2	1.13	0.90
4:D:60:ARG:HH11	4:D:66:ASP:HA	1.38	0.89
4:D:99:ARG:HA	4:D:99:ARG:HE	1.39	0.88
4:D:12:VAL:HG12	4:D:110:GLU:HB2	1.59	0.85

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	19/21 (90%)	11 (58%)	6 (32%)	2 (10%)	0 7
2	B	19/30 (63%)	16 (84%)	2 (10%)	1 (5%)	2 17
3	C	116/118 (98%)	108 (93%)	3 (3%)	5 (4%)	2 22
4	D	112/114 (98%)	104 (93%)	6 (5%)	2 (2%)	8 41
5	E	282/317 (89%)	261 (93%)	19 (7%)	2 (1%)	22 61

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	F	13/16 (81%)	11 (85%)	1 (8%)	1 (8%)	1	10
All	All	561/616 (91%)	511 (91%)	37 (7%)	13 (2%)	6	36

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
2	B	8	GLY
3	C	100	TYR
3	C	105	MET
4	D	33	SER

### 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	20/20 (100%)	15 (75%)	5 (25%)	0	4
2	B	17/26 (65%)	15 (88%)	2 (12%)	5	25
3	C	101/101 (100%)	97 (96%)	4 (4%)	31	64
4	D	100/100 (100%)	92 (92%)	8 (8%)	12	41
5	E	268/293 (92%)	253 (94%)	15 (6%)	21	54
6	F	15/16 (94%)	9 (60%)	6 (40%)	0	1
All	All	521/556 (94%)	481 (92%)	40 (8%)	13	42

5 of 40 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	E	194	THR
6	F	707	ASP
5	E	211	GLU
5	E	250	ASP
6	F	712	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
5	E	123	ASN
5	E	152	ASN
5	E	281	ASN
5	E	247	HIS
4	D	43	GLN

### 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](#)

8 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	$\# Z  > 2$
7	NAG	G	1	5,7	14,14,15	0.33	0	17,19,21	1.22	2 (11%)
7	NAG	G	2	7	14,14,15	0.44	0	17,19,21	0.62	0
7	BMA	G	3	7	11,11,12	0.21	0	15,15,17	0.73	0
7	MAN	G	4	7	11,11,12	0.32	0	15,15,17	0.60	0
8	NAG	H	1	8,5	14,14,15	0.35	0	17,19,21	1.15	2 (11%)
8	NAG	H	2	8	14,14,15	0.22	0	17,19,21	0.93	1 (5%)
8	BMA	H	3	8	11,11,12	0.30	0	15,15,17	0.67	0
8	FUC	H	4	8	10,10,11	0.43	0	14,14,16	0.64	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
7	NAG	G	1	5,7	-	0/6/23/26	0/1/1/1
7	NAG	G	2	7	-	0/6/23/26	0/1/1/1
7	BMA	G	3	7	-	2/2/19/22	0/1/1/1
7	MAN	G	4	7	-	0/2/19/22	0/1/1/1
8	NAG	H	1	8,5	-	0/6/23/26	0/1/1/1
8	NAG	H	2	8	-	0/6/23/26	0/1/1/1
8	BMA	H	3	8	-	0/2/19/22	0/1/1/1
8	FUC	H	4	8	-	-	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	G	1	NAG	C1-O5-C5	3.23	116.57	112.19
8	H	1	NAG	C1-C2-N2	-3.03	105.31	110.49
8	H	2	NAG	O5-C1-C2	-2.77	106.91	111.29
8	H	1	NAG	O5-C1-C2	2.18	114.73	111.29
7	G	1	NAG	O4-C4-C3	2.05	115.09	110.35

There are no chirality outliers.

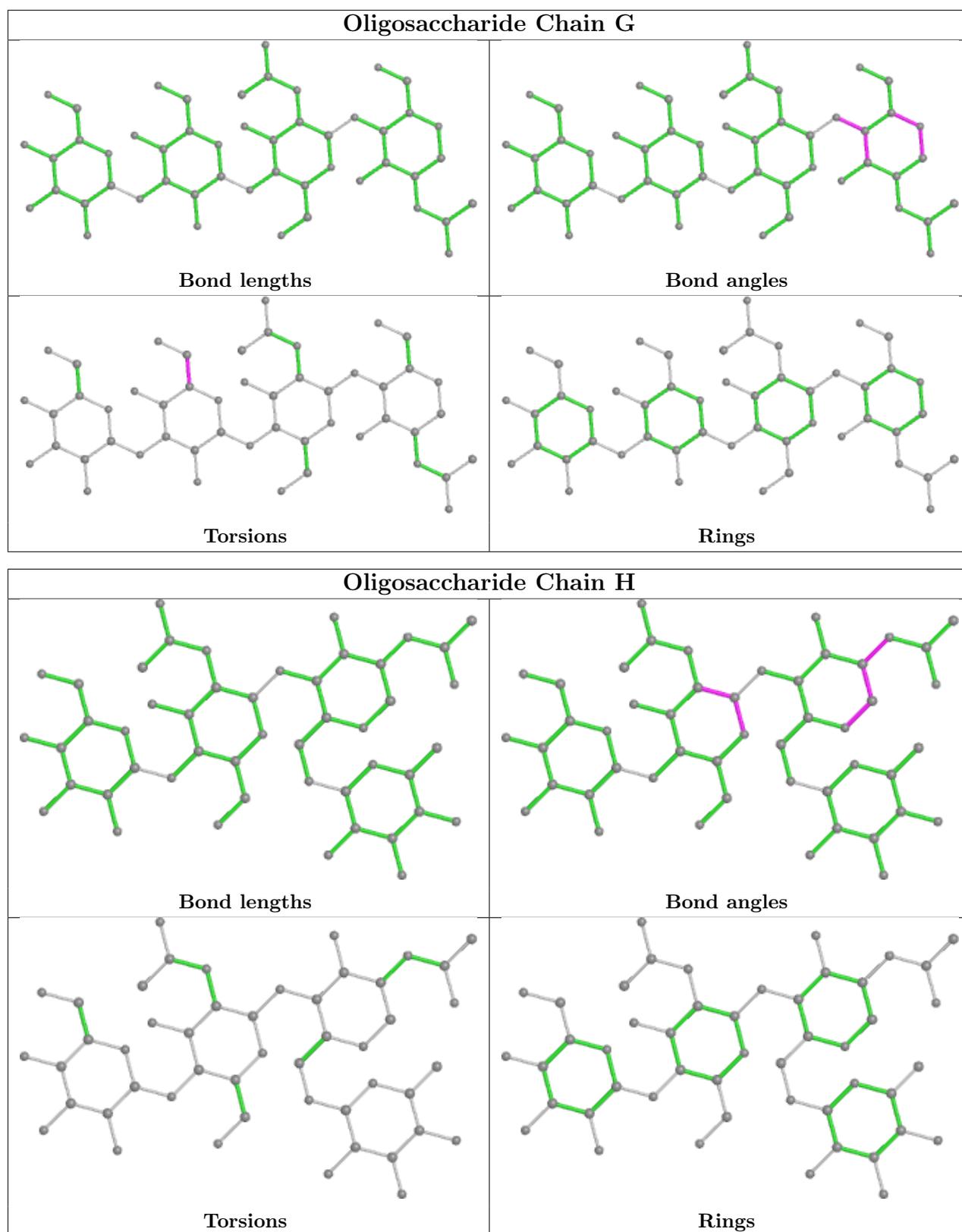
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	G	3	BMA	C4-C5-C6-O6
7	G	3	BMA	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 [i](#)

3 ligands 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	$\# Z  > 2$
9	NAG	E	507	5	14,14,15	0.37	0	17,19,21	1.29	2 (11%)
9	NAG	E	502	5	14,14,15	0.31	0	17,19,21	0.79	1 (5%)
9	NAG	E	501	5	14,14,15	0.35	0	17,19,21	0.84	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	E	507	5	-	1/6/23/26	0/1/1/1
9	NAG	E	502	5	-	1/6/23/26	0/1/1/1
9	NAG	E	501	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	E	507	NAG	O5-C1-C2	-4.07	104.86	111.29
9	E	501	NAG	C1-O5-C5	3.12	116.41	112.19
9	E	507	NAG	C2-N2-C7	2.48	126.44	122.90
9	E	502	NAG	C1-O5-C5	2.37	115.40	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	E	502	NAG	O5-C5-C6-O6
9	E	507	NAG	C3-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [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	21/21 (100%)	-0.27	1 (4%) 30 27	221, 244, 251, 260	0
2	B	21/30 (70%)	-0.50	1 (4%) 30 27	214, 232, 258, 277	0
3	C	118/118 (100%)	0.22	9 (7%) 13 14	195, 266, 299, 304	0
4	D	114/114 (100%)	-0.01	6 (5%) 26 24	198, 230, 265, 276	0
5	E	288/317 (90%)	-0.23	3 (1%) 82 77	183, 225, 279, 300	0
6	F	15/16 (93%)	-0.12	0 100 100	196, 219, 265, 277	0
All	All	577/616 (93%)	-0.10	20 (3%) 44 39	183, 235, 289, 304	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	117	SER	6.8
3	C	96	ALA	5.5
4	D	40	ALA	5.2
3	C	35	ASN	4.7
3	C	95	CYS	4.2

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

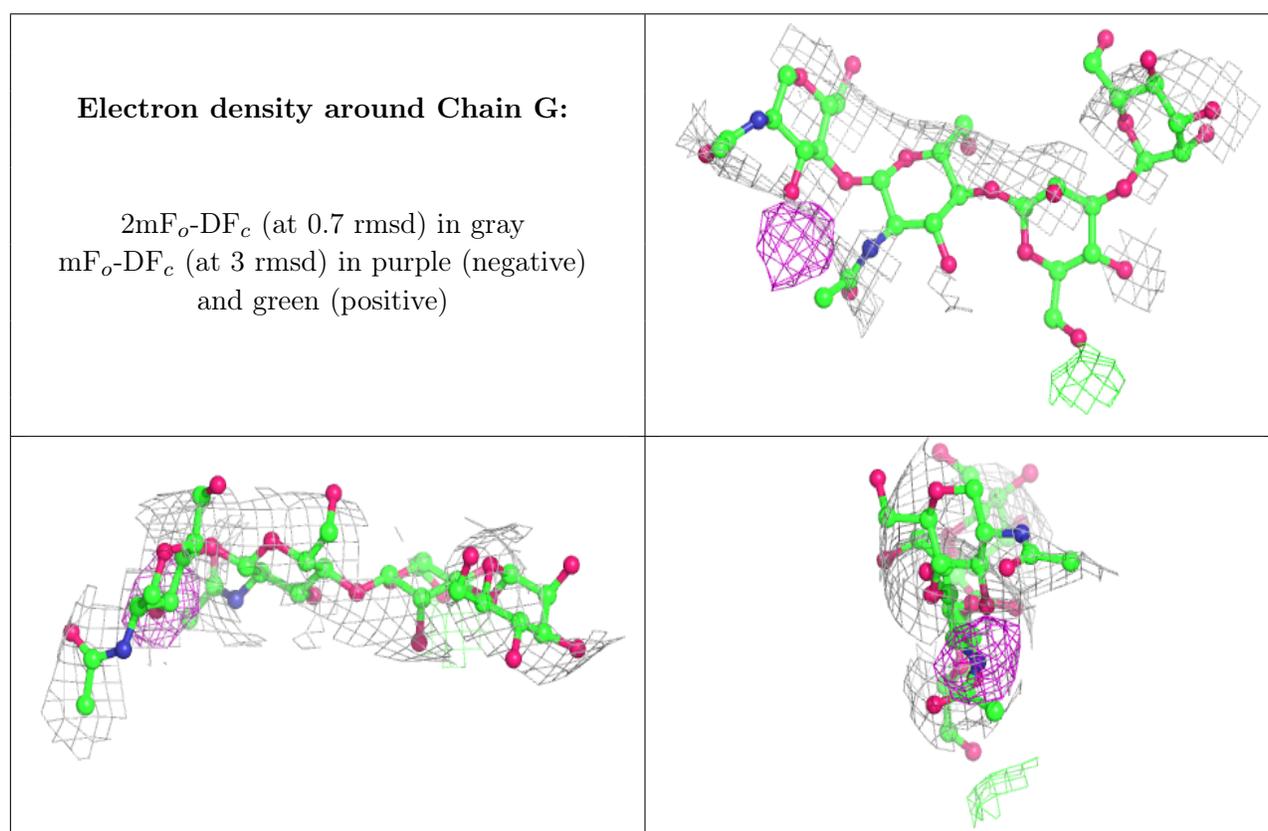
There are no non-standard protein/DNA/RNA residues in this entry.

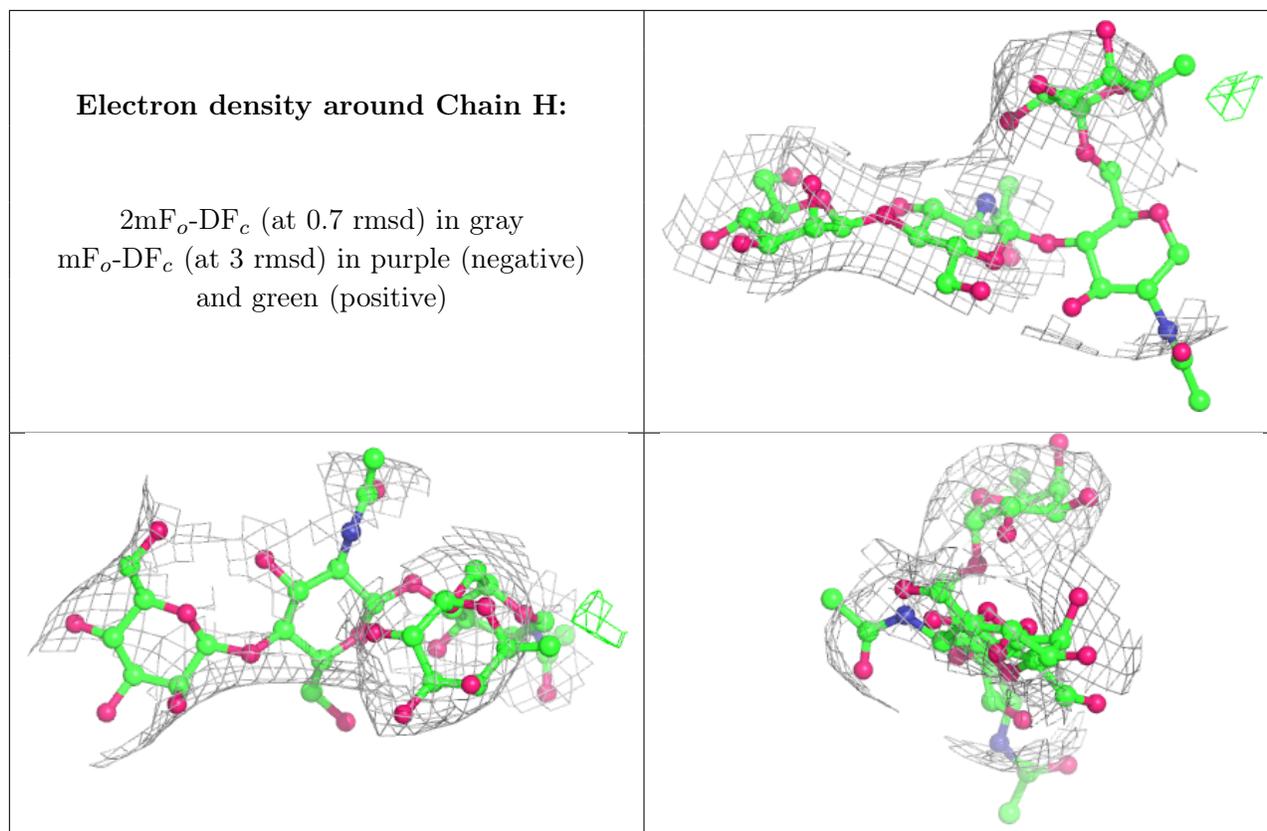
### 6.3 Carbohydrates [i](#)

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( $\text{\AA}^2$ )	$Q < 0.9$
8	BMA	H	3	11/12	0.68	0.44	310,315,316,316	0
7	NAG	G	1	14/15	0.84	0.21	211,215,221,224	0
7	BMA	G	3	11/12	0.86	0.30	250,256,261,263	0
8	NAG	H	1	14/15	0.87	0.20	263,274,279,288	0
8	NAG	H	2	14/15	0.88	0.38	294,297,303,307	0
7	MAN	G	4	11/12	0.90	0.49	264,266,270,273	0
8	FUC	H	4	10/11	0.93	0.20	259,265,268,270	0
7	NAG	G	2	14/15	0.94	0.24	228,234,239,243	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.





## 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(Å <sup>2</sup> )	Q<0.9
9	NAG	E	502	14/15	0.81	0.22	282,293,297,297	0
9	NAG	E	501	14/15	0.83	0.32	294,297,299,300	0
9	NAG	E	507	14/15	0.88	0.21	244,256,262,265	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.