Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 08:30 AM BST

PDB ID	:	6U3M
Title	:	DQ2-P.fluor-alpha1a
Authors	:	Petersen, J.; Rossjohn, J.
Deposited on	:	2019-08-22
Resolution	:	1.90 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ⓘ) were used in the production of this report:

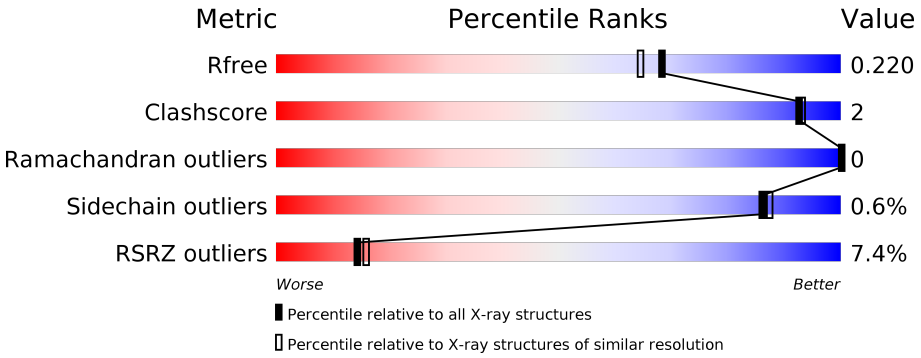
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 >=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 <=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	191	3% 92% • 5%
1	C	191	% 92% • 5%
2	B	206	11% 82% 6% 12%
2	D	206	11% 83% • 12%
3	E	21	5% 52% 48%
3	F	21	5% 52% 48%

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Mol	Chain	Length	Quality of chain
4	G	2	100%
4	H	2	100%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6586 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 HLA class II histocompatibility antigen, DQ alpha 1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	181	Total	C	N	O	S	0	0	0
			1445	931	236	276	2			
1	A	181	Total	C	N	O	S	0	0	0
			1445	931	236	276	2			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	44	SER	CYS	conflict	UNP P01909
C	182	THR	-	expression tag	UNP P01909
C	183	SER	-	expression tag	UNP P01909
C	184	GLY	-	expression tag	UNP P01909
C	185	ASP	-	expression tag	UNP P01909
C	186	ASP	-	expression tag	UNP P01909
C	187	ASP	-	expression tag	UNP P01909
C	188	ASP	-	expression tag	UNP P01909
C	189	LYS	-	expression tag	UNP P01909
A	44	SER	CYS	conflict	UNP P01909
A	182	THR	-	expression tag	UNP P01909
A	183	SER	-	expression tag	UNP P01909
A	184	GLY	-	expression tag	UNP P01909
A	185	ASP	-	expression tag	UNP P01909
A	186	ASP	-	expression tag	UNP P01909
A	187	ASP	-	expression tag	UNP P01909
A	188	ASP	-	expression tag	UNP P01909
A	189	LYS	-	expression tag	UNP P01909

- Molecule 2 is a protein called MHC class II HLA-DQ-beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	181	Total	C	N	O	S	0	0	0
			1474	932	261	274	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	181	Total	C	N	O	S	0	0	0
			1474	932	261	274	7			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	GLY	-	expression tag	UNP O19712
D	-4	GLY	-	expression tag	UNP O19712
D	-3	SER	-	expression tag	UNP O19712
D	-2	GLY	-	expression tag	UNP O19712
D	-1	ALA	-	expression tag	UNP O19712
D	0	SER	-	expression tag	UNP O19712
D	193	THR	-	expression tag	UNP O19712
D	194	GLY	-	expression tag	UNP O19712
D	195	GLY	-	expression tag	UNP O19712
D	196	ASP	-	expression tag	UNP O19712
D	197	ASP	-	expression tag	UNP O19712
D	198	ASP	-	expression tag	UNP O19712
D	199	ASP	-	expression tag	UNP O19712
D	200	LYS	-	expression tag	UNP O19712
B	-5	GLY	-	expression tag	UNP O19712
B	-4	GLY	-	expression tag	UNP O19712
B	-3	SER	-	expression tag	UNP O19712
B	-2	GLY	-	expression tag	UNP O19712
B	-1	ALA	-	expression tag	UNP O19712
B	0	SER	-	expression tag	UNP O19712
B	193	THR	-	expression tag	UNP O19712
B	194	GLY	-	expression tag	UNP O19712
B	195	GLY	-	expression tag	UNP O19712
B	196	ASP	-	expression tag	UNP O19712
B	197	ASP	-	expression tag	UNP O19712
B	198	ASP	-	expression tag	UNP O19712
B	199	ASP	-	expression tag	UNP O19712
B	200	LYS	-	expression tag	UNP O19712

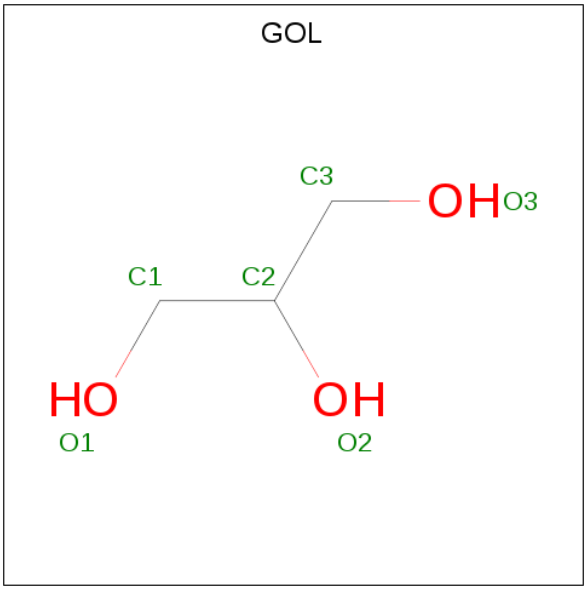
- Molecule 3 is a protein called Alpha1a peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	11	Total	C	N	O	S	0	0	0
			85	58	11	14	2			
3	F	11	Total	C	N	O	S	0	0	0
			85	58	11	14	2			

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

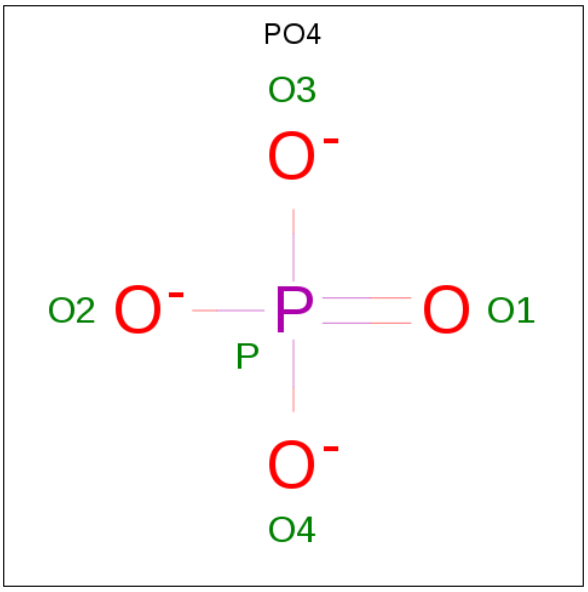
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
4	G	2	Total C N O 28 16 2 10	0	0	0
4	H	2	Total C N O 28 16 2 10	0	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



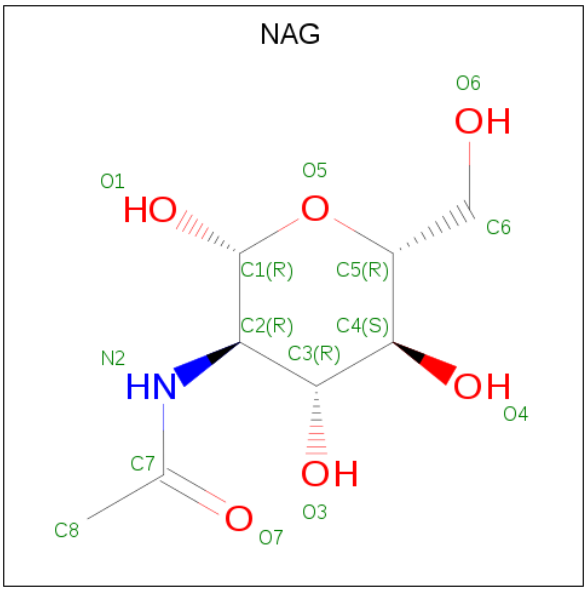
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total C O 6 3 3	0	0
5	C	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
6	D	1	5	4	1	0	0

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
7	A	1	14	8	1	5	0	0

- Molecule 8 is water.

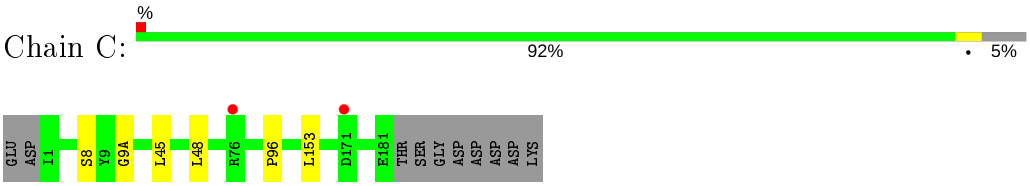
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	C	149	Total O 149 149	0	0
8	D	111	Total O 111 111	0	0
8	A	130	Total O 130 130	0	0
8	B	77	Total O 77 77	0	0
8	E	7	Total O 7 7	0	0
8	F	5	Total O 5 5	0	0

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.

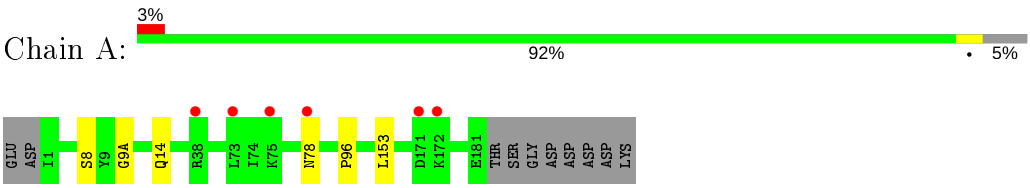
- Molecule 1: HLA class II histocompatibility antigen, DQ alpha 1 chain

Chain C:



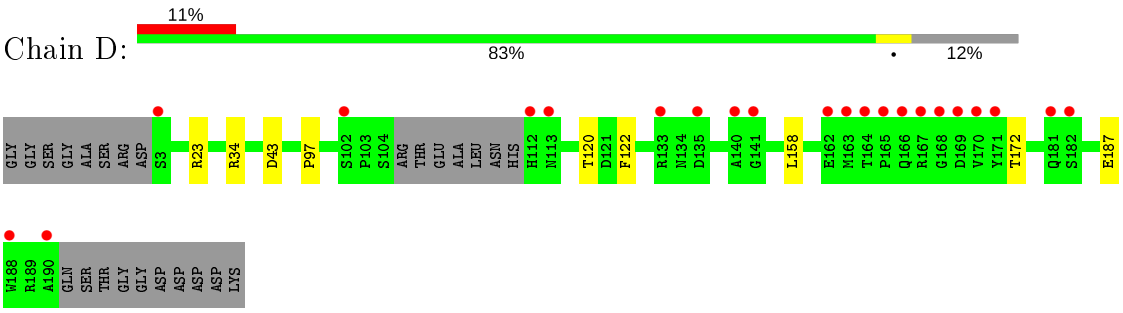
- Molecule 1: HLA class II histocompatibility antigen, DQ alpha 1 chain

Chain A:



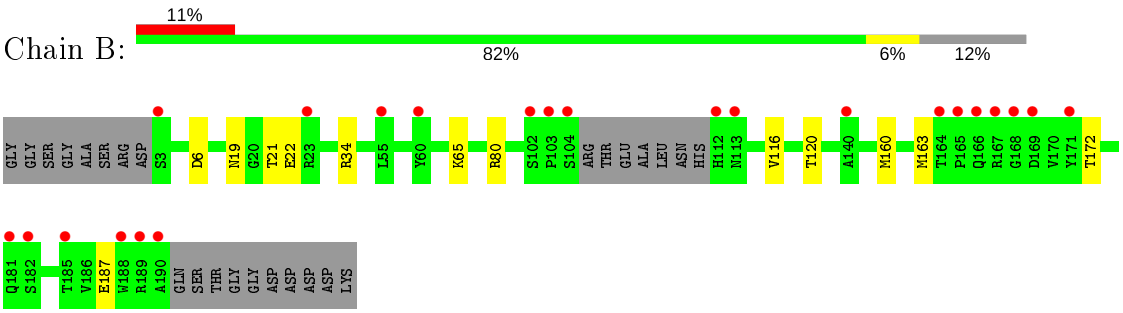
- Molecule 2: MHC class II HLA-DQ-beta-1

Chain D:



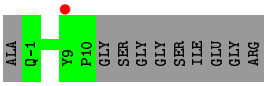
- Molecule 2: MHC class II HLA-DQ-beta-1

Chain B:



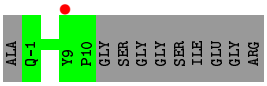
- Molecule 3: Alpha1a peptide

Chain E:



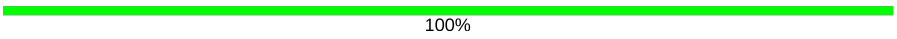
• Molecule 3: Alpha1a peptide

Chain F:



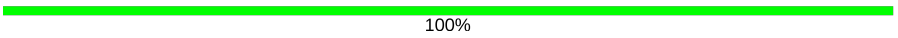
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:



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

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	94.94Å 96.28Å 105.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.47 – 1.90 48.14 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.47-1.90) 99.9 (48.14-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.72 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.190 , 0.220 0.190 , 0.220	Depositor DCC
R_{free} test set	1800 reflections (2.34%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 47.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6586	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.42 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.2980e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, PO4, NAG

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	0.37	0/1487	0.55	0/2031
1	C	0.42	0/1487	0.57	0/2031
2	B	0.34	0/1507	0.54	0/2049
2	D	0.38	0/1507	0.57	0/2049
3	E	0.32	0/90	0.47	0/125
3	F	0.34	0/90	0.48	0/125
All	All	0.38	0/6168	0.56	0/8410

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1445	0	1397	4	0
1	C	1445	0	1397	3	0
2	B	1474	0	1432	8	0
2	D	1474	0	1431	5	0
3	E	85	0	83	0	0
3	F	85	0	83	0	0
4	G	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	28	0	25	0	0
5	A	12	0	15	0	0
5	C	12	0	15	1	0
6	D	5	0	0	0	0
7	A	14	0	13	0	0
8	A	130	0	0	1	0
8	B	77	0	0	1	0
8	C	149	0	0	1	0
8	D	111	0	0	1	0
8	E	7	0	0	0	0
8	F	5	0	0	0	0
All	All	6586	0	5916	18	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 (18) 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:19:ASN:ND2	2:B:22:GLU:OE1	2.14	0.80
2:D:172:THR:HG22	2:D:187:GLU:HG2	1.75	0.68
2:B:172:THR:HG22	2:B:187:GLU:HG2	1.77	0.67
2:B:21:THR:O	2:B:80:ARG:NH1	2.27	0.64
5:C:1101:GOL:O3	8:C:1201:HOH:O	2.17	0.55
2:B:116:VAL:HG22	2:B:160:MET:HG2	1.89	0.53
1:A:96:PRO:HD3	2:B:120:THR:HG21	1.93	0.51
1:A:78:ASN:ND2	8:A:1104:HOH:O	2.46	0.47
2:B:65:LYS:HD2	2:B:65:LYS:HA	1.66	0.44
2:D:23:ARG:NH2	2:D:43:ASP:OD2	2.48	0.44
1:C:96:PRO:HD3	2:D:120:THR:HG21	1.99	0.43
1:A:14:GLN:NE2	2:B:6:ASP:OD2	2.52	0.43
2:D:34:ARG:NH1	8:D:1103:HOH:O	2.41	0.43
1:A:8:SER:C	1:A:9(A):GLY:HA2	2.40	0.42
2:B:34:ARG:NH1	8:B:306:HOH:O	2.47	0.42
1:C:8:SER:C	1:C:9(A):GLY:HA2	2.41	0.41
1:C:45:LEU:HB3	1:C:48:LEU:HG	2.03	0.41
2:D:97:PRO:HB3	2:D:122:PHE:HB3	2.02	0.41

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	179/191 (94%)	177 (99%)	2 (1%)	0	100	100
1	C	179/191 (94%)	176 (98%)	3 (2%)	0	100	100
2	B	177/206 (86%)	171 (97%)	6 (3%)	0	100	100
2	D	177/206 (86%)	171 (97%)	6 (3%)	0	100	100
3	E	9/21 (43%)	8 (89%)	1 (11%)	0	100	100
3	F	9/21 (43%)	8 (89%)	1 (11%)	0	100	100
All	All	730/836 (87%)	711 (97%)	19 (3%)	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	165/174 (95%)	164 (99%)	1 (1%)	86	87
1	C	165/174 (95%)	164 (99%)	1 (1%)	86	87
2	B	163/184 (89%)	162 (99%)	1 (1%)	86	87
2	D	163/184 (89%)	162 (99%)	1 (1%)	86	87
3	E	10/16 (62%)	10 (100%)	0	100	100
3	F	10/16 (62%)	10 (100%)	0	100	100
All	All	676/748 (90%)	672 (99%)	4 (1%)	86	87

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

Mol	Chain	Res	Type
1	C	153	LEU
2	D	158	LEU
1	A	153	LEU
2	B	163	MET

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

Mol	Chain	Res	Type
1	A	68	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

4 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	$\#\|Z\| > 2$
4	NAG	G	1	1,4	14,14,15	0.34	0	17,19,21	0.49	0
4	NAG	G	2	4	14,14,15	0.42	0	17,19,21	0.44	0
4	NAG	H	1	2,4	14,14,15	0.57	0	17,19,21	0.42	0
4	NAG	H	2	4	14,14,15	0.39	0	17,19,21	0.44	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
4	NAG	G	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	NAG	H	1	2,4	-	1/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

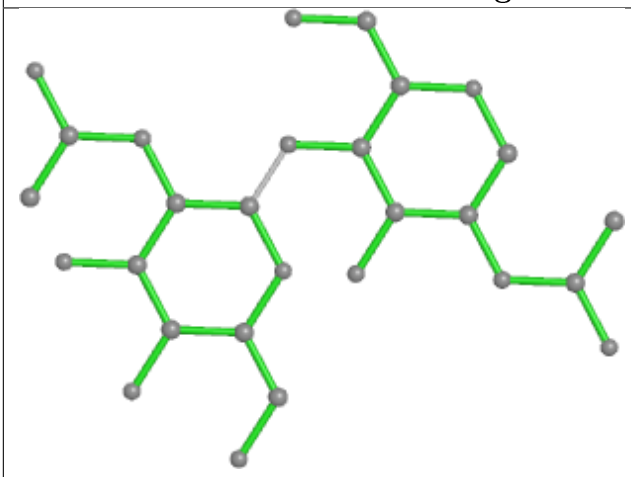
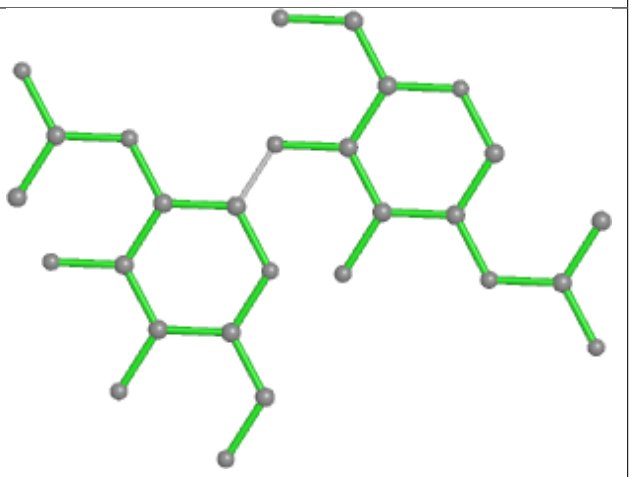
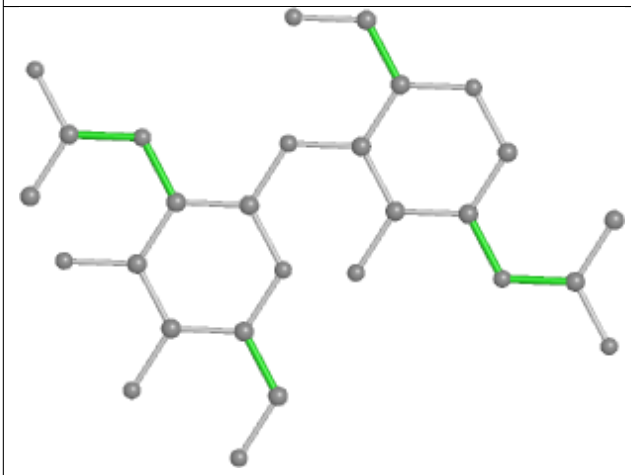
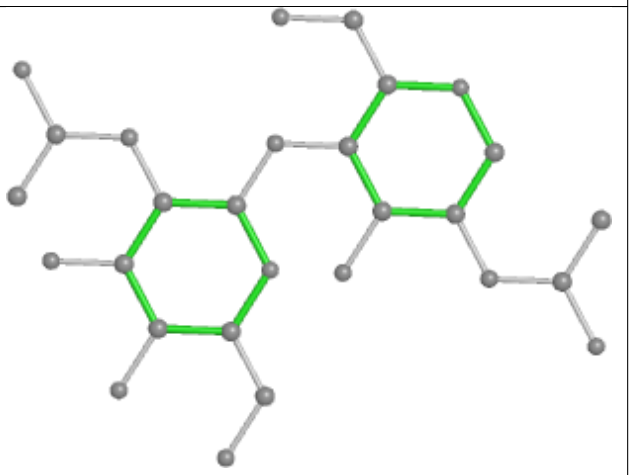
All (3) torsion outliers are listed below:

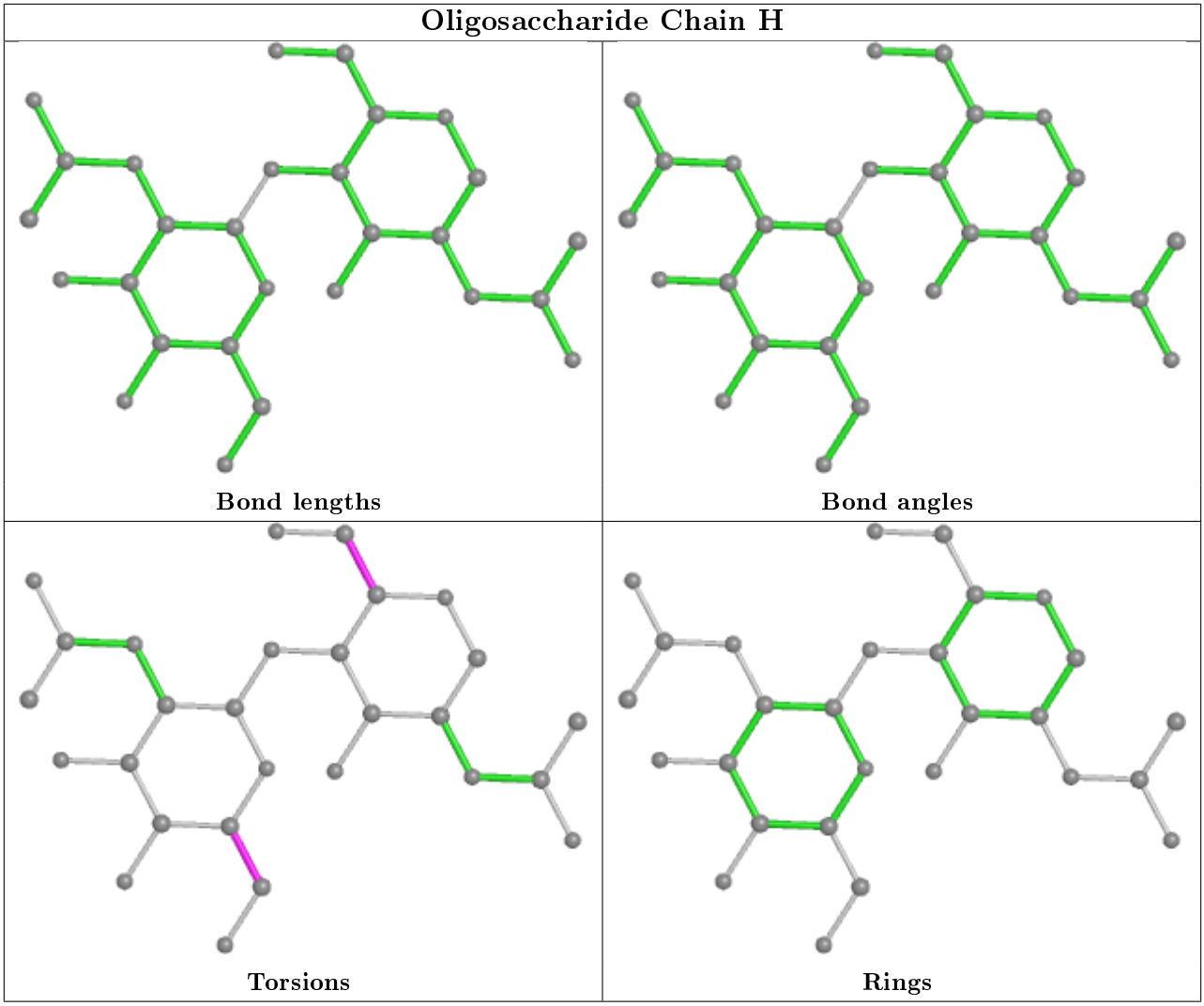
Mol	Chain	Res	Type	Atoms
4	H	2	NAG	C4-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6

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.

Oligosaccharide Chain G	
 Bond lengths	 Bond angles
 Torsions	 Rings



5.6 Ligand geometry (i)

6 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	$\#\|Z\| > 2$
6	PO4	D	1003	-	4,4,4	0.82	0	6,6,6	0.47	0
5	GOL	A	1003	-	5,5,5	1.01	0	5,5,5	1.48	1 (20%)
5	GOL	A	1002	-	5,5,5	1.35	1 (20%)	5,5,5	1.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
7	NAG	A	1001	1	14,14,15	0.40	0	17,19,21	0.44	0
5	GOL	C	1101	-	5,5,5	1.29	1 (20%)	5,5,5	1.02	0
5	GOL	C	1102	-	5,5,5	1.11	0	5,5,5	1.21	1 (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
5	GOL	A	1003	-	-	4/4/4/4	-
7	NAG	A	1001	1	-	2/6/23/26	0/1/1/1
5	GOL	A	1002	-	-	2/4/4/4	-
5	GOL	C	1101	-	-	2/4/4/4	-
5	GOL	C	1102	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1002	GOL	O2-C2	-2.58	1.35	1.43
5	C	1101	GOL	O2-C2	-2.45	1.36	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1003	GOL	C3-C2-C1	-2.54	101.82	111.70
5	C	1102	GOL	C3-C2-C1	-2.05	103.74	111.70

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1101	GOL	O1-C1-C2-C3
5	A	1003	GOL	O1-C1-C2-O2
7	A	1001	NAG	O5-C5-C6-O6
5	A	1003	GOL	O1-C1-C2-C3
5	A	1003	GOL	C1-C2-C3-O3
7	A	1001	NAG	C4-C5-C6-O6
5	A	1003	GOL	O2-C2-C3-O3
5	C	1101	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	A	1002	GOL	O1-C1-C2-C3
5	A	1002	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	1101	GOL	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^{th} 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^2)	Q<0.9
1	A	181/191 (94%)	0.16	6 (3%) 46 49	26, 43, 73, 94	0
1	C	181/191 (94%)	0.19	2 (1%) 80 82	22, 35, 59, 83	0
2	B	181/206 (87%)	0.65	23 (12%) 3 4	24, 55, 87, 119	0
2	D	181/206 (87%)	0.72	22 (12%) 4 4	22, 45, 88, 134	0
3	E	11/21 (52%)	0.59	1 (9%) 9 10	33, 45, 86, 87	0
3	F	11/21 (52%)	0.45	1 (9%) 9 10	28, 37, 78, 83	0
All	All	746/836 (89%)	0.43	55 (7%) 14 16	22, 44, 83, 134	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	112	HIS	16.4
2	B	112	HIS	11.2
2	D	167	ARG	7.9
2	D	164	THR	6.7
2	B	167	ARG	6.2
3	E	9	TYR	6.0
2	D	190	ALA	5.2
2	D	113	ASN	5.1
2	B	168	GLY	5.0
2	D	163	MET	4.9
2	B	164	THR	4.9
2	D	165	PRO	4.6
2	B	190	ALA	4.6
2	B	140	ALA	4.5
2	B	113	ASN	4.4
2	D	166	GLN	4.4
2	B	165	PRO	4.1
1	A	73	LEU	4.1
2	B	3	SER	4.1

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Mol	Chain	Res	Type	RSRZ
2	D	171	TYR	4.0
2	D	168	GLY	3.8
2	D	141	GLY	3.7
2	B	166	GLN	3.6
3	F	9	TYR	3.4
2	D	181	GLN	3.2
1	C	171	ASP	3.2
2	D	188	TRP	3.1
1	A	78	ASN	3.1
2	D	140	ALA	3.1
1	A	171	ASP	3.1
2	D	182	SER	2.9
2	D	133	ARG	2.9
2	D	170	VAL	2.9
2	B	181	GLN	2.9
2	D	3	SER	2.9
2	B	188	TRP	2.8
2	D	102	SER	2.7
2	B	103	PRO	2.7
1	A	172	LYS	2.7
2	B	169	ASP	2.6
1	C	76	ARG	2.4
2	B	23	ARG	2.4
2	B	182	SER	2.4
2	B	189	ARG	2.3
2	D	169	ASP	2.3
2	B	55	LEU	2.2
1	A	38	ARG	2.2
2	B	60	TYR	2.2
2	B	185	THR	2.2
1	A	75	LYS	2.2
2	B	171	TYR	2.1
2	B	104	SER	2.1
2	D	135	ASP	2.1
2	B	102	SER	2.0
2	D	162	GLU	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)

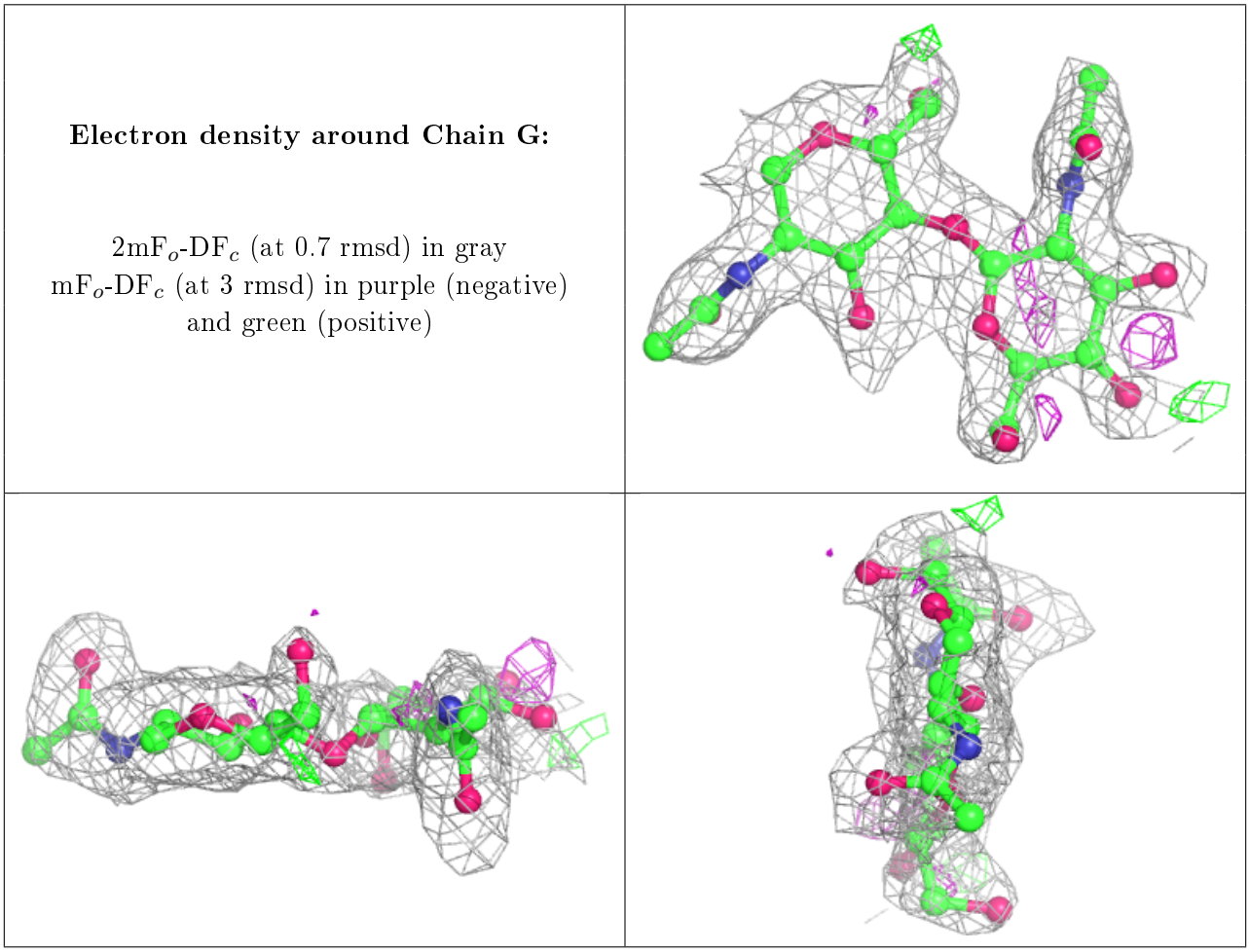
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NAG	G	2	14/15	0.61	0.28	72,75,80,81	0
4	NAG	H	1	14/15	0.67	0.21	49,61,66,73	0
4	NAG	H	2	14/15	0.68	0.33	76,79,81,82	0
4	NAG	G	1	14/15	0.89	0.14	42,48,59,63	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.

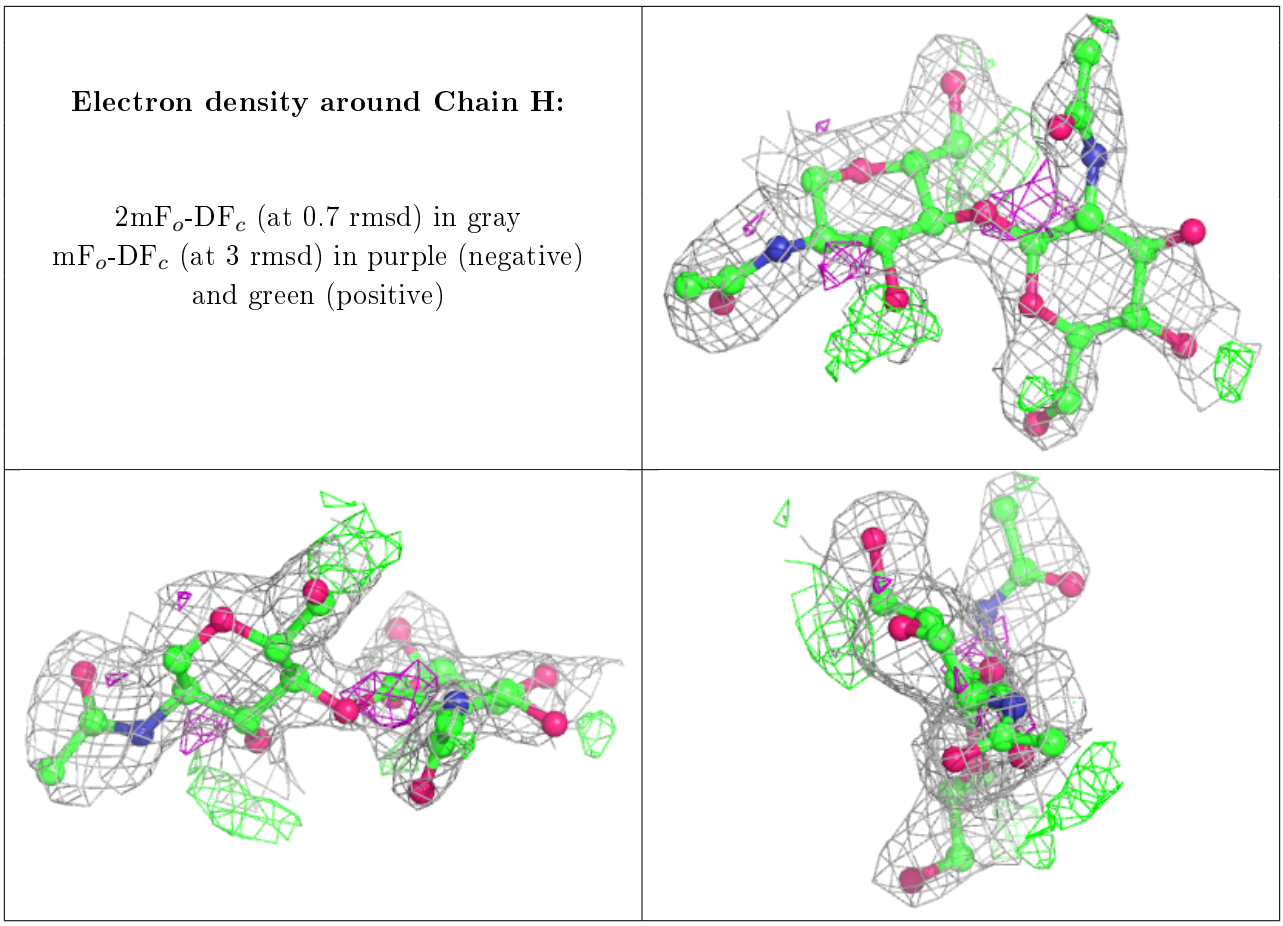
Electron density around Chain G:

$2mF_o$-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)



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^{th} 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^2)	Q<0.9
5	GOL	A	1003	6/6	0.76	0.34	70,74,77,81	0
5	GOL	C	1102	6/6	0.80	0.35	67,77,79,81	0
7	NAG	A	1001	14/15	0.83	0.20	47,55,63,63	0
5	GOL	C	1101	6/6	0.86	0.34	57,64,69,73	0
5	GOL	A	1002	6/6	0.88	0.26	58,60,61,65	0
6	PO4	D	1003	5/5	0.97	0.10	51,51,53,55	0

6.5 Other polymers

There are no such residues in this entry.