



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 12:50 AM UTC

PDB ID : 5Y3K / pdb_00005y3k
Title : Crystal structure of horse TLR9 in complex with two DNAs (CpG DNA and GCGCAC DNA)
Authors : Ohto, U.; Shimizu, T.
Deposited on : 2017-07-29
Resolution : 2.70 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

A user guide is available at

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

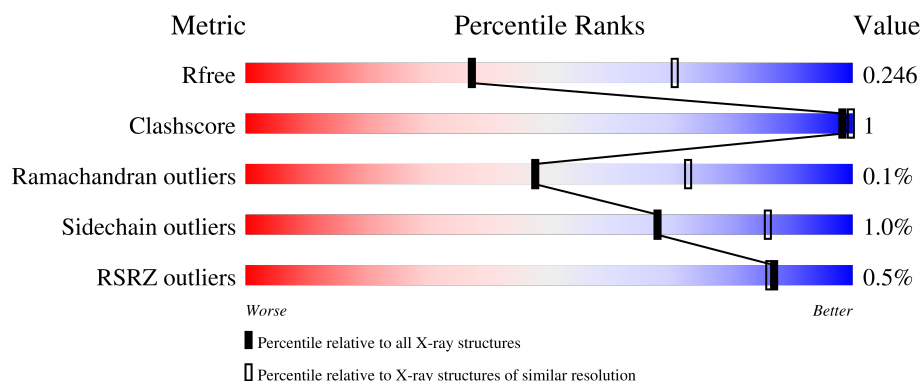
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 ≥ 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	802	
1	B	802	
2	C	10	
2	D	10	
3	E	6	

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Mol	Chain	Length	Quality of chain
3	F	6	 100%
4	G	2	 50% 50%
4	H	2	 100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12766 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 Toll-like receptor 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	750	Total	C	N	O	S	0	0	0
			5893	3765	1038	1062	28			
1	B	750	Total	C	N	O	S	0	0	0
			5893	3765	1038	1062	28			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ARG	-	expression tag	UNP Q2EEY0
A	23	SER	-	expression tag	UNP Q2EEY0
A	24	PRO	-	expression tag	UNP Q2EEY0
A	25	TRP	-	expression tag	UNP Q2EEY0
A	818	GLU	CYS	conflict	UNP Q2EEY0
A	820	LEU	-	expression tag	UNP Q2EEY0
A	821	VAL	-	expression tag	UNP Q2EEY0
A	822	PRO	-	expression tag	UNP Q2EEY0
A	823	ARG	-	expression tag	UNP Q2EEY0
B	22	ARG	-	expression tag	UNP Q2EEY0
B	23	SER	-	expression tag	UNP Q2EEY0
B	24	PRO	-	expression tag	UNP Q2EEY0
B	25	TRP	-	expression tag	UNP Q2EEY0
B	818	GLU	CYS	conflict	UNP Q2EEY0
B	820	LEU	-	expression tag	UNP Q2EEY0
B	821	VAL	-	expression tag	UNP Q2EEY0
B	822	PRO	-	expression tag	UNP Q2EEY0
B	823	ARG	-	expression tag	UNP Q2EEY0

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*GP*GP*CP*GP*TP*TP*TP*TP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	P	0	0	0
			203	99	33	62	9			
2	D	10	Total	C	N	O	P	0	0	0
			203	99	33	62	9			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*CP*GP*CP*AP*C)-3').

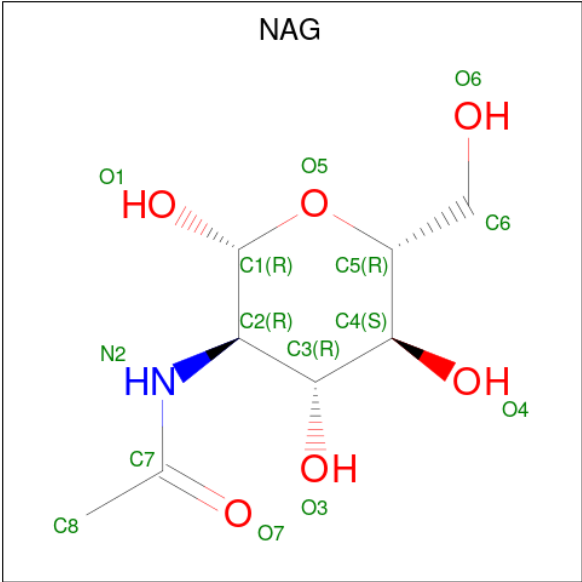
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	6	Total	C	N	O	P	0	0	0
			119	57	24	33	5			
3	F	6	Total	C	N	O	P	0	0	0
			119	57	24	33	5			

- 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	0	0	0
			28	16	2	10			
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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

- Molecule 6 is water.

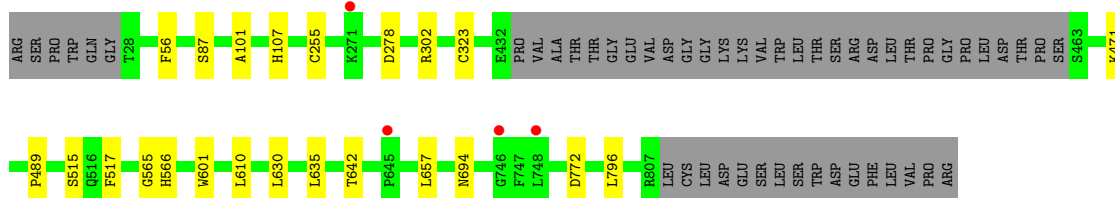
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	35	Total	O	0	0
			35	35		
6	E	1	Total	O	0	0
			1	1		
6	B	33	Total	O	0	0
			33	33		
6	F	1	Total	O	0	0
			1	1		

3 Residue-property plots

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.

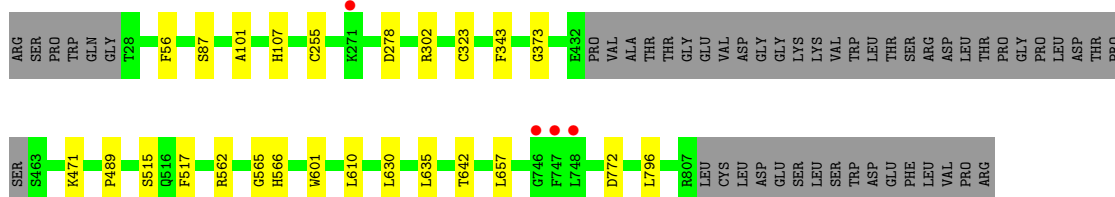
- Molecule 1: Toll-like receptor 9

Chain A: 



- Molecule 1: Toll-like receptor 9

Chain B: 

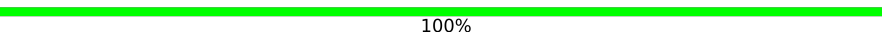


- Molecule 2: DNA (5'-D(*AP*GP*GP*CP*GP*TP*TP*TP*TP*T)-3')

Chain C: 

There are no outlier residues recorded for this chain.

- Molecule 2: DNA (5'-D(*AP*GP*GP*CP*GP*TP*TP*TP*TP*T)-3')

Chain D: 

There are no outlier residues recorded for this chain.

- Molecule 3: DNA (5'-D(*GP*CP*GP*CP*AP*C)-3')

Chain E: 

There are no outlier residues recorded for this chain.

- Molecule 3: DNA (5'-D(*GP*CP*GP*CP*AP*C)-3')

Chain F:  100%

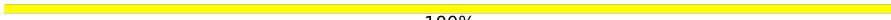
There are no outlier residues recorded for this chain.

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

Chain G:  50% 50%

 NAG1
NAG2

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

Chain H:  100%

 NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.85Å 126.50Å 140.72Å 90.00° 98.49° 90.00°	Depositor
Resolution (Å)	47.14 – 2.70 47.14 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.8 (47.14-2.70) 98.8 (47.14-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.205 , 0.246 0.209 , 0.246	Depositor DCC
R_{free} test set	2533 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	54.3	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12766	wwPDB-VP
Average B, all atoms (Å ²)	64.0	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 47.10 % 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 1.0408e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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	0.70	0/6032	0.85	3/8206 (0.0%)
1	B	0.69	0/6032	0.85	3/8206 (0.0%)
2	C	0.44	0/226	0.94	0/348
2	D	0.44	0/226	0.86	0/348
3	E	0.42	0/133	0.70	0/203
3	F	0.46	0/133	0.76	0/203
All	All	0.68	0/12782	0.85	6/17514 (0.0%)

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	517	PHE	CA-C-N	5.33	123.60	120.24
1	A	517	PHE	C-N-CA	5.33	123.60	120.24
1	B	517	PHE	CA-C-N	5.15	123.48	120.24
1	B	517	PHE	C-N-CA	5.15	123.48	120.24
1	B	471	LYS	N-CA-C	5.09	118.27	111.75

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	5893	0	5916	8	0
1	B	5893	0	5915	8	0
2	C	203	0	117	0	0
2	D	203	0	117	0	0
3	E	119	0	68	0	0
3	F	119	0	68	0	0
4	G	28	0	25	0	0
4	H	28	0	25	0	0
5	A	98	0	91	1	0
5	B	112	0	104	0	0
6	A	35	0	0	0	0
6	B	33	0	0	0	0
6	E	1	0	0	0	0
6	F	1	0	0	0	0
All	All	12766	0	12446	16	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 1.

The worst 5 of 16 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:A:489:PRO:HG3	1:A:515:SER:O	2.08	0.54
1:B:489:PRO:HG3	1:B:515:SER:O	2.12	0.48
1:B:278:ASP:OD1	1:B:302:ARG:NH1	2.51	0.44
1:A:278:ASP:OD1	1:A:302:ARG:NH1	2.51	0.44
1:B:657:LEU:HD23	1:B:657:LEU:C	2.43	0.43

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	746/802 (93%)	701 (94%)	44 (6%)	1 (0%)	48	73
1	B	746/802 (93%)	700 (94%)	45 (6%)	1 (0%)	48	73
All	All	1492/1604 (93%)	1401 (94%)	89 (6%)	2 (0%)	48	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	PHE
1	B	56	PHE

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	666/712 (94%)	660 (99%)	6 (1%)	70	87
1	B	666/712 (94%)	659 (99%)	7 (1%)	65	85
All	All	1332/1424 (94%)	1319 (99%)	13 (1%)	68	86

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	255	CYS
1	B	323	CYS
1	B	796	LEU
1	B	642	THR
1	B	772	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	154	ASN
1	B	381	GLN
1	B	650	ASN
1	B	335	GLN

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Mol	Chain	Res	Type
1	B	385	GLN

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	0.51	0	17,19,21	1.37	2 (11%)
4	NAG	G	2	4	14,14,15	0.54	0	17,19,21	1.03	0
4	NAG	H	1	1,4	14,14,15	0.51	0	17,19,21	1.33	3 (17%)
4	NAG	H	2	4	14,14,15	0.66	0	17,19,21	1.21	3 (17%)

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	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	3/6/23/26	0/1/1/1
4	NAG	H	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1	NAG	C1-O5-C5	2.84	115.99	112.19
4	H	2	NAG	O3-C3-C2	2.63	114.86	109.40
4	H	1	NAG	C1-C2-N2	-2.55	106.42	110.43
4	G	1	NAG	C2-N2-C7	2.51	126.27	122.90
4	H	2	NAG	C2-N2-C7	2.24	125.91	122.90

There are no chirality outliers.

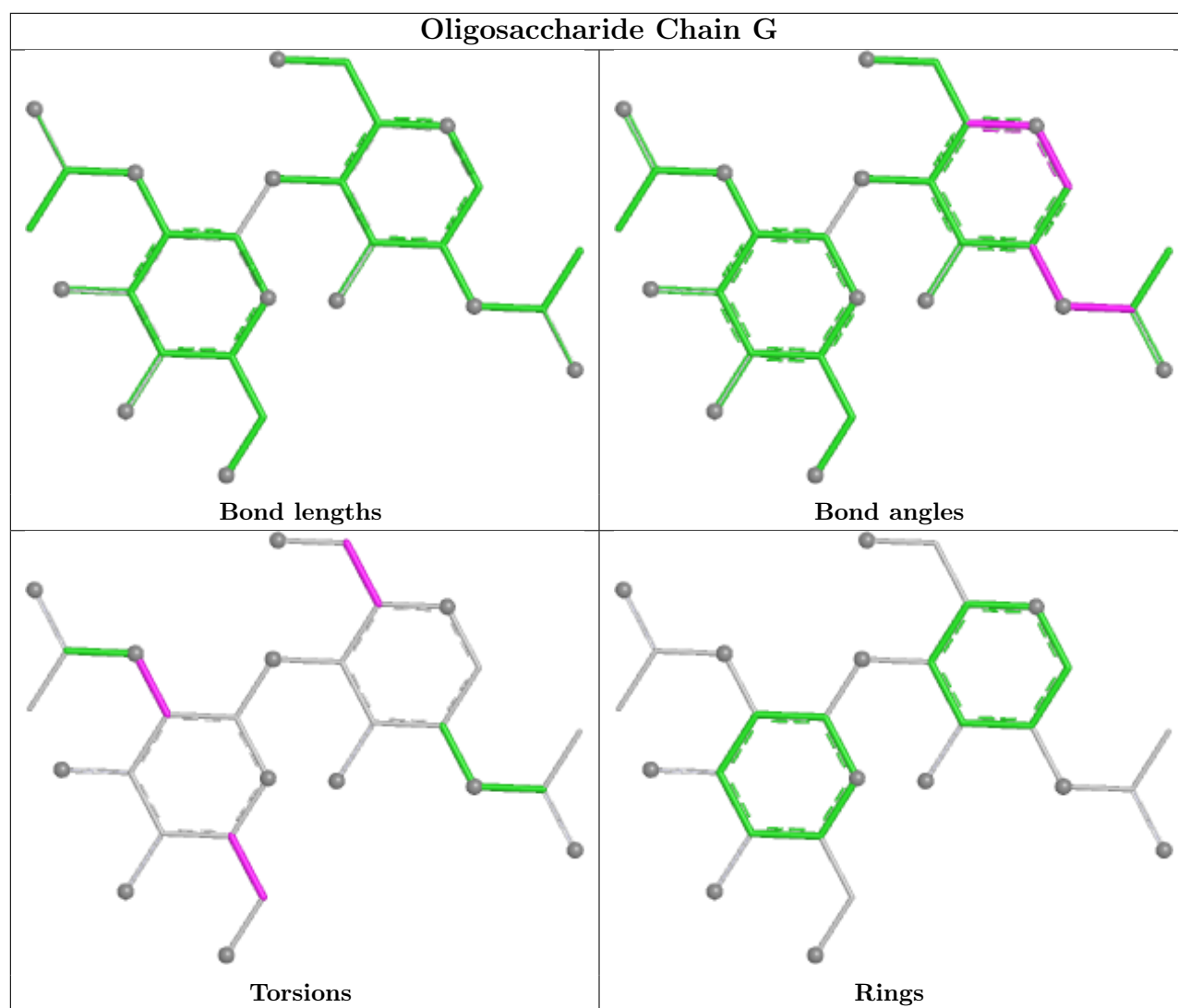
5 of 9 torsion outliers are listed below:

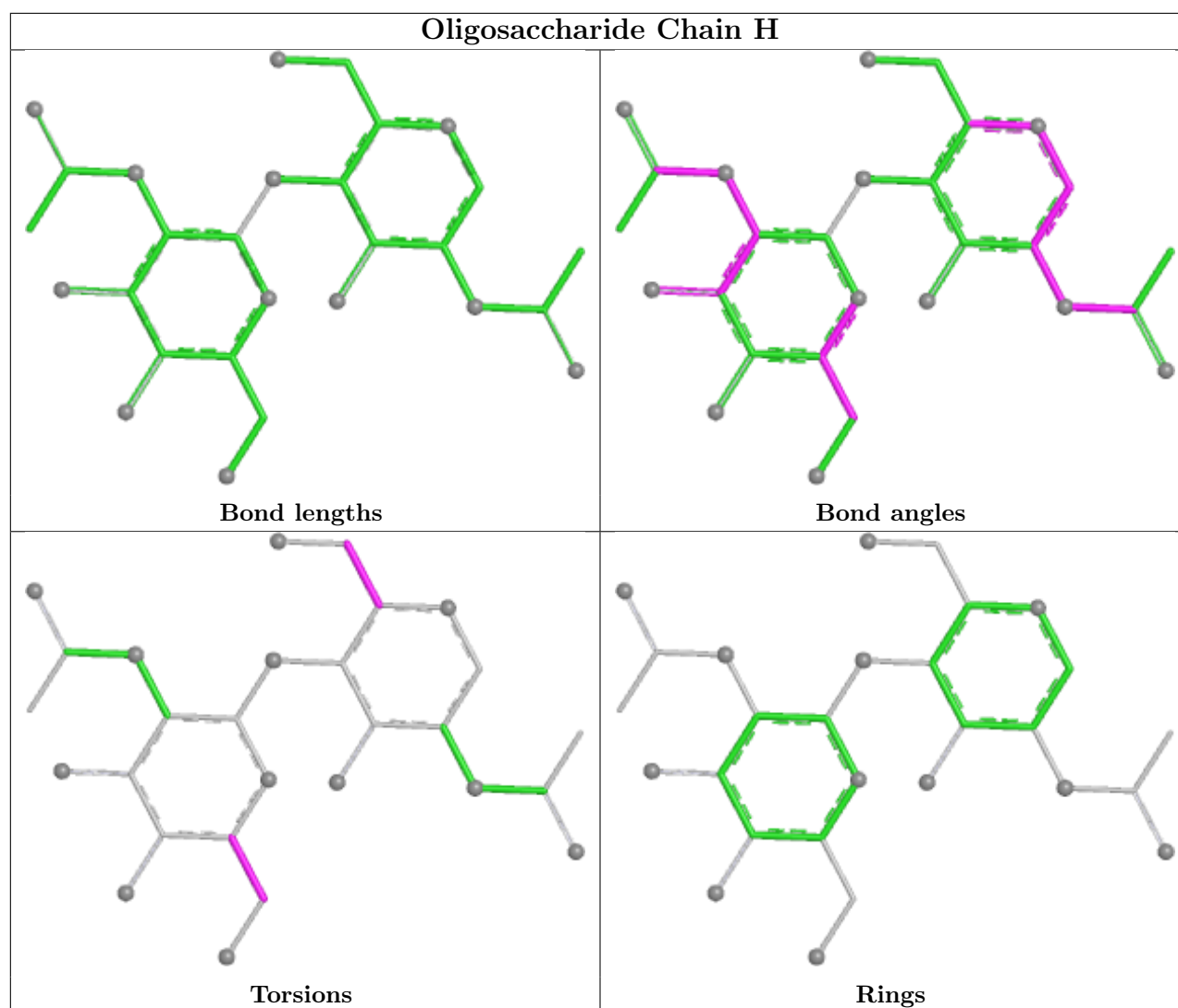
Mol	Chain	Res	Type	Atoms
4	H	1	NAG	O5-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6

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 [i](#)

15 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	$\# Z > 2$
5	NAG	A	1004	1	14,14,15	0.54	0	17,19,21	2.11	4 (23%)
5	NAG	A	1008	1	14,14,15	0.45	0	17,19,21	1.13	0
5	NAG	A	1006	1	14,14,15	0.49	0	17,19,21	1.15	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	B	1006	1	14,14,15	0.76	1 (7%)	17,19,21	1.04	0
5	NAG	A	1005	1	14,14,15	0.37	0	17,19,21	0.85	0
5	NAG	A	1009	1	14,14,15	0.31	0	17,19,21	1.08	1 (5%)
5	NAG	B	1005	1	14,14,15	0.66	0	17,19,21	1.38	2 (11%)
5	NAG	A	1007	1	14,14,15	0.53	0	17,19,21	1.48	3 (17%)
5	NAG	B	1010	1	14,14,15	0.78	0	17,19,21	1.73	2 (11%)
5	NAG	A	1003	1	14,14,15	0.52	0	17,19,21	1.75	5 (29%)
5	NAG	B	1009	1	14,14,15	0.52	0	17,19,21	1.16	2 (11%)
5	NAG	B	1008	1	14,14,15	0.32	0	17,19,21	1.01	1 (5%)
5	NAG	B	1004	1	14,14,15	0.43	0	17,19,21	1.73	3 (17%)
5	NAG	B	1007	1	14,14,15	0.54	0	17,19,21	1.47	3 (17%)
5	NAG	B	1003	1	14,14,15	0.47	0	17,19,21	1.45	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
5	NAG	A	1004	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1008	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1006	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1006	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1005	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1009	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1005	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1007	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1010	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1003	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1009	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1008	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1004	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1007	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1003	1	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1006	NAG	C1-C2	2.05	1.55	1.52

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1004	NAG	C1-O5-C5	6.70	121.17	112.19
5	B	1004	NAG	C1-O5-C5	4.82	118.64	112.19
5	B	1010	NAG	C2-N2-C7	4.34	128.72	122.90
5	A	1003	NAG	O5-C1-C2	-4.01	105.08	111.29
5	B	1007	NAG	C1-C2-N2	-3.99	104.15	110.43

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1006	NAG	O5-C5-C6-O6
5	B	1006	NAG	C4-C5-C6-O6
5	A	1004	NAG	O5-C5-C6-O6
5	B	1004	NAG	C4-C5-C6-O6
5	B	1004	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1009	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 [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, 95th 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	750/802 (93%)	-0.33	4 (0%) 87 86	33, 59, 88, 111	0
1	B	750/802 (93%)	-0.26	4 (0%) 87 86	35, 64, 93, 114	0
2	C	10/10 (100%)	-0.38	0 100 100	67, 74, 105, 106	0
2	D	10/10 (100%)	-0.29	0 100 100	66, 71, 100, 102	0
3	E	6/6 (100%)	-0.65	0 100 100	45, 47, 73, 105	0
3	F	6/6 (100%)	-0.20	0 100 100	51, 55, 81, 123	0
All	All	1532/1636 (93%)	-0.30	8 (0%) 87 86	33, 62, 92, 123	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	746	GLY	3.1
1	A	645	PRO	2.6
1	A	748	LEU	2.5
1	B	748	LEU	2.3
1	B	747	PHE	2.3

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

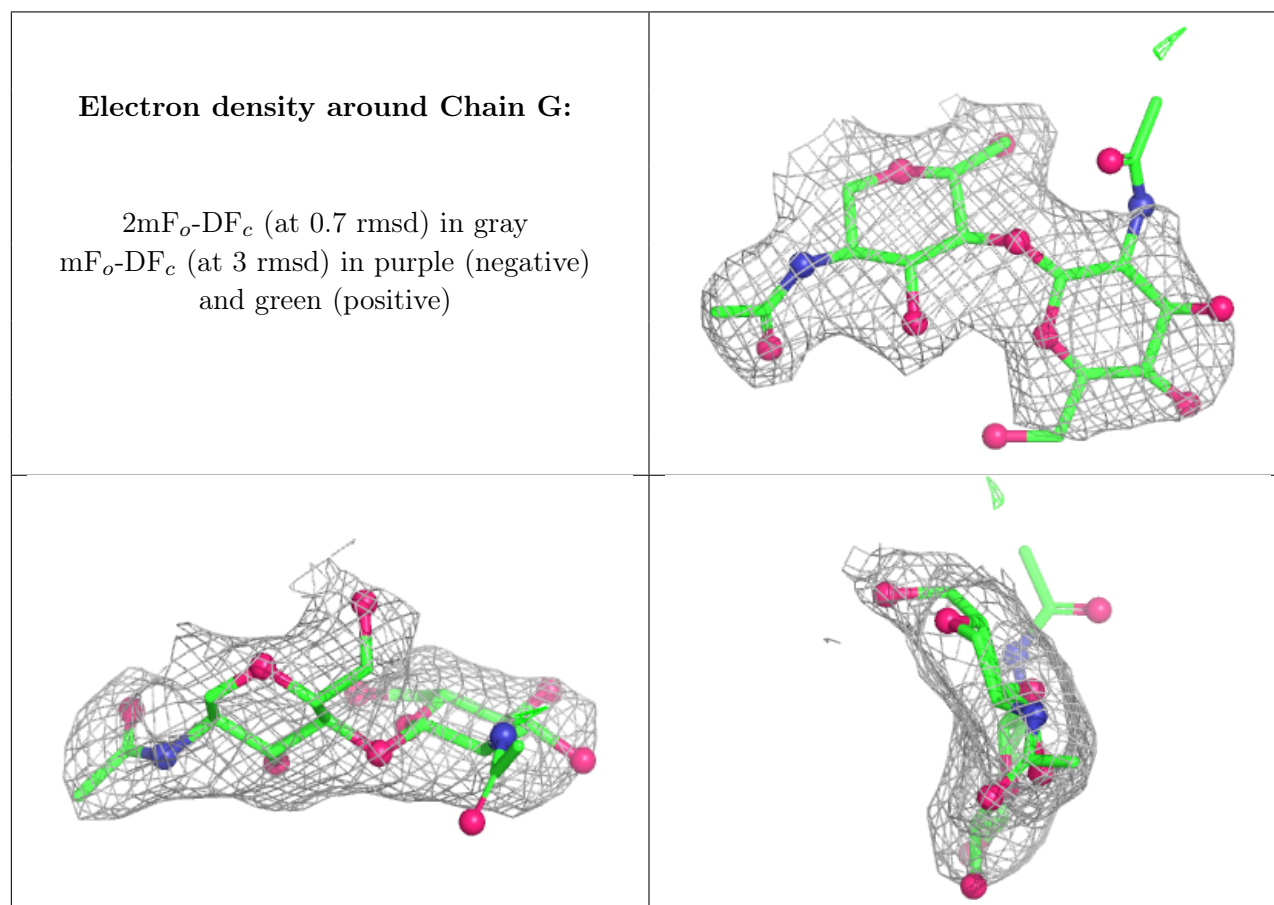
There are no non-standard protein/DNA/RNA residues in this entry.

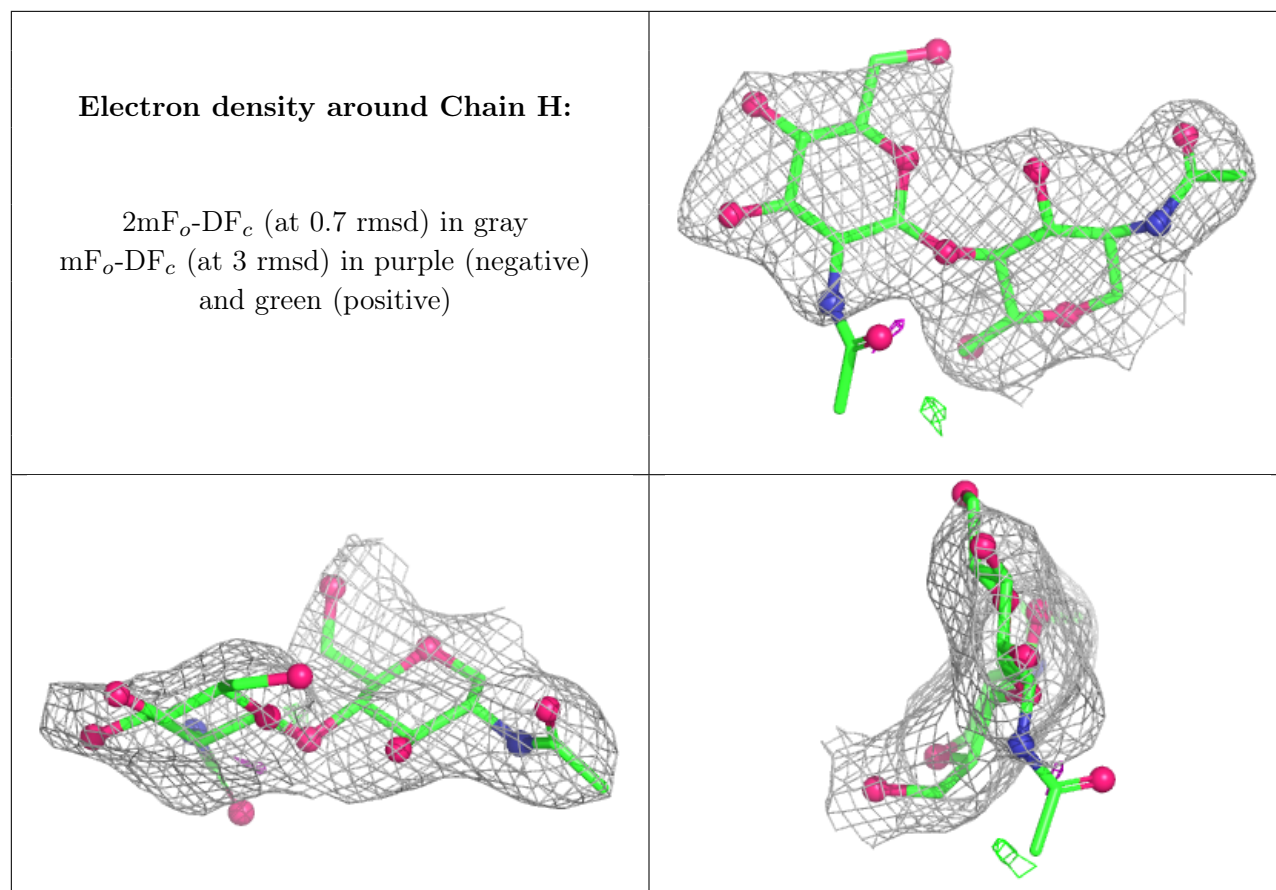
6.3 Carbohydrates [i](#)

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, 95th 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	2	14/15	0.81	0.11	80,85,99,101	0
4	NAG	G	2	14/15	0.85	0.10	81,89,99,99	0
4	NAG	G	1	14/15	0.93	0.06	59,64,67,71	0
4	NAG	H	1	14/15	0.94	0.06	58,63,68,76	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 ⓘ

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, 95th 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	NAG	B	1010	14/15	0.74	0.13	119,127,130,132	0
5	NAG	B	1005	14/15	0.79	0.10	86,98,102,106	0
5	NAG	A	1003	14/15	0.80	0.10	76,83,85,85	0
5	NAG	A	1005	14/15	0.83	0.09	84,94,96,97	0
5	NAG	B	1007	14/15	0.84	0.10	74,78,82,84	0
5	NAG	A	1007	14/15	0.86	0.10	71,75,79,81	0
5	NAG	A	1006	14/15	0.87	0.10	78,83,85,87	0
5	NAG	B	1006	14/15	0.87	0.10	86,91,93,94	0
5	NAG	B	1003	14/15	0.89	0.08	66,73,76,77	0
5	NAG	B	1004	14/15	0.90	0.08	72,78,84,88	0
5	NAG	B	1008	14/15	0.91	0.08	60,64,66,70	0
5	NAG	A	1008	14/15	0.92	0.09	55,57,59,60	0
5	NAG	A	1004	14/15	0.92	0.07	72,76,80,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	B	1009	14/15	0.94	0.07	52,56,62,63	0
5	NAG	A	1009	14/15	0.95	0.06	50,54,56,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.