

Full wwPDB X-ray Structure Validation Report (i)

Dec 20, 2023 – 01:01 PM EST

PDB ID : 4MDI

Title : Immune Receptor

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Deposited on : 2013-08-22

Resolution : 2.00 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 (i) 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 (1)) 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 : 2.36

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)

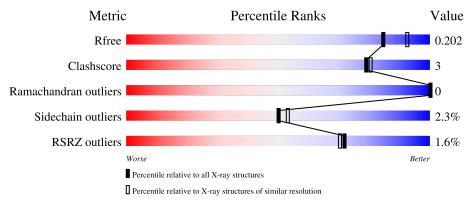
Validation Pipeline (wwPDB-VP) : 2.36

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 2.00 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{\rm A})}) \end{array}$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	189	86%	8% • 5%
2	В	200	87%	6% • 6%
3	С	13	85%	15%
4	D	2	100%	



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 3672 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, DR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total 1542	C 997	N 247	O 293	S 5	0	10	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	THR	-	expression tag	UNP P01903
A	183	SER	-	expression tag	UNP P01903
A	184	GLY	-	expression tag	UNP P01903
A	185	ASP	-	expression tag	UNP P01903
A	186	ASP	-	expression tag	UNP P01903
A	187	ASP	-	expression tag	UNP P01903
A	188	ASP	-	expression tag	UNP P01903
A	189	LYS	-	expression tag	UNP P01903

• Molecule 2 is a protein called HLA class II histocompatibility antigen, DRB1-4 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	В	189	Total 1577	C 995	N 277	O 300	S 5	0	1	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	-1	GLY	-	expression tag	UNP P13760
В	0	SER	- expression to		UNP P13760
В	67	ILE	LEU	variant	UNP P13760
В	70	ASP	GLN	variant	UNP P13760
В	71	GLU	LYS	variant	UNP P13760
В	86	VAL	GLY	variant	UNP P13760
В	191	THR	-	expression tag	UNP P13760
В	192	GLY	-	expression tag	UNP P13760



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Chain	Residue	Modelled	Actual	Comment	Reference
В	193	GLY	-	expression tag	UNP P13760
В	194	ASP	-	expression tag	UNP P13760
В	195	ASP	-	expression tag	UNP P13760
В	196	ASP	-	expression tag	UNP P13760
В	197	ASP	-	expression tag	UNP P13760
В	198	LYS	-	expression tag	UNP P13760

• Molecule 3 is a protein called Citrullinated Vimentin.

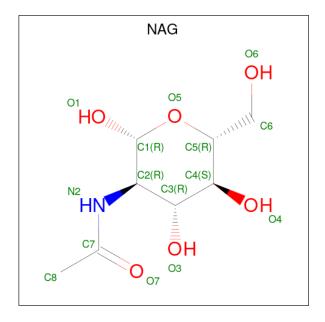
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	С	13	Total 96	C 58	N 21	O 17	0	0	0

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



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

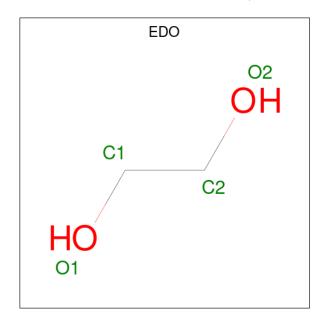
• Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).





Mol	Chain	Residues	A	ton	\mathbf{ns}		ZeroOcc	AltConf
5	A	1	Total 14			O 5	0	0
5	В	1	Total 14	C 8		O 5	0	0

 \bullet Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\mathrm{C_2H_6O_2}).$



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	В	1	Total C 4 2	O 2	0	0

• Molecule 7 is water.

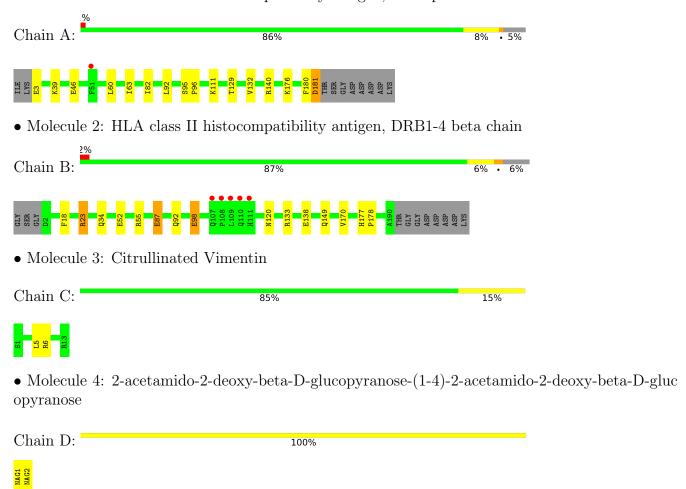
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	199	Total O 199 199	0	0
7	В	189	Total O 189 189	0	0
7	С	9	Total O 9 9	0	0



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, DR alpha chain





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants	67.02Å 182.93Å 77.38Å	Donogitor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.37 - 2.00	Depositor
Resolution (A)	39.37 - 2.00	EDS
% Data completeness	100.0 (39.37-2.00)	Depositor
(in resolution range)	100.0 (39.37-2.00)	EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.97 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
D D	0.161 , 0.203	Depositor
R, R_{free}	0.163 , 0.202	DCC
R_{free} test set	1684 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 54.2	EDS
L-test for twinning ²	$< L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3672	wwPDB-VP
Average B, all atoms $(Å^2)$	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: CIR, EDO, 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		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.36	0/1605	0.52	0/2186	
2	В	0.39	0/1621	0.54	0/2206	
3	С	0.36	0/84	0.60	0/111	
All	All	0.37	0/3310	0.53	0/4503	

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1542	0	1476	9	0
2	В	1577	0	1468	11	0
3	С	96	0	103	0	0
4	D	28	0	25	0	0
5	A	14	0	13	0	0
5	В	14	0	13	0	0
6	В	4	0	6	0	0
7	A	199	0	0	0	0
7	В	189	0	0	1	0
7	С	9	0	0	0	0
All	All	3672	0	3104	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 3.

All (18) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} (\mathring{\rm A}) \end{array}$	Clash overlap (Å)
1:A:46[B]:GLU:H	1:A:46[B]:GLU:CD	2.00	0.63
1:A:180:PHE:O	1:A:181:ASP:HB2	2.01	0.60
2:B:98[A]:GLU:CG	2:B:120:ASN:HB2	2.32	0.60
1:A:39:LYS:HG2	1:A:60:LEU:HD21	1.90	0.54
2:B:149:GLN:NE2	7:B:414:HOH:O	2.26	0.48
2:B:52:GLU:HA	2:B:55:ARG:HE	1.81	0.45
2:B:98[A]:GLU:HG2	2:B:120:ASN:HB2	1.99	0.44
1:A:95:SER:HB2	1:A:96:PRO:HD2	1.99	0.43
1:A:176:LYS:HD3	1:A:176:LYS:HA	1.89	0.43
2:B:18:PHE:HB2	2:B:23:ARG:HB3	2.00	0.43
1:A:3:GLU:HA	2:B:18:PHE:CD1	2.55	0.42
1:A:111:LYS:HG2	1:A:140:ARG:CZ	2.49	0.42
2:B:87:GLU:OE1	2:B:92:GLN:NE2	2.53	0.41
2:B:133:ARG:NH1	2:B:138:GLU:OE2	2.49	0.41
1:A:82:ILE:HD11	2:B:34:GLN:HE22	1.86	0.41
2:B:177:HIS:CD2	2:B:178:PRO:HD2	2.56	0.41
2:B:98[A]:GLU:HG3	2:B:120:ASN:HB2	2.03	0.40
1:A:129:THR:O	1:A:132:VAL:HG22	2.22	0.40

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	Analysed Favoured Allowed		Outliers	Perce	ntiles
1	A	187/189 (99%)	185 (99%)	2 (1%)	0	100	100
2	В	188/200 (94%)	183 (97%)	5 (3%)	0	100	100
3	С	10/13 (77%)	10 (100%)	0	0	100	100



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Mol	Chain	Analysed Favoured Allowed		Outliers	Perce	ntiles	
All	All	385/402 (96%)	378 (98%)	7 (2%)	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	174/173 (101%)	171 (98%)	3 (2%)	60 65		
2	В	172/179 (96%)	167 (97%)	5 (3%)	42 43		
3	С	10/10 (100%)	9 (90%)	1 (10%)	7 4		
All	All	356/362 (98%)	347 (98%)	9 (2%)	50 49		

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ILE
1	A	92	LEU
1	A	181	ASP
2	В	23	ARG
2	В	87	GLU
2	В	98[A]	GLU
2	В	98[B]	GLU
2	В	170	VAL
3	С	5	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
2	В	34	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)

1 non-standard protein/DNA/RNA residue 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	B	ond leng	$_{ m gths}$	В	ond ang	gles
WIOI	Type	Chain	rtes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CIR	С	6	3	9,10,11	3.20	3 (33%)	6,11,13	1.88	3 (50%)

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	CIR	С	6	3	-	3/8/9/11	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(\mathbf{\mathring{A}})$	$\operatorname{Ideal}(\text{\AA})$
3	С	6	CIR	C7-N6	7.66	1.44	1.34
3	С	6	CIR	C7-N8	4.74	1.44	1.33
3	С	6	CIR	С3-СА	-2.35	1.50	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
3	С	6	CIR	C5-N6-C7	-2.91	119.27	122.73
3	С	6	CIR	N8-C7-N6	2.44	119.43	116.85
3	С	6	CIR	O7-C7-N8	-2.20	119.44	123.22

There are no chirality outliers.



All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	С	6	CIR	O7-C7-N6-C5
3	С	6	CIR	N8-C7-N6-C5
3	С	6	CIR	C3-C4-C5-N6

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

2 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	Dag	Dec	Pog	Pos	Dog	Dog	Ros	Link	Bo	ond leng	$ ag{ths}$	В	ond ang	cles
	туре			LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2							
4	NAG	D	1	4,1	14,14,15	0.54	0	17,19,21	1.70	4 (23%)							
4	NAG	D	2	4	14,14,15	0.72	0	17,19,21	1.74	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	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mo	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	D	1	NAG	O4-C4-C3	-4.26	100.51	110.35
4	D	2	NAG	C3-C4-C5	3.91	117.22	110.24



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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	D	2	NAG	C4-C3-C2	3.65	116.37	111.02
4	D	2	NAG	O5-C1-C2	-3.64	105.53	111.29
4	D	1	NAG	C1-O5-C5	3.12	116.42	112.19
4	D	1	NAG	C4-C3-C2	2.71	114.99	111.02
4	D	1	NAG	C3-C4-C5	2.32	114.39	110.24

There are no chirality outliers.

All (4) torsion outliers are listed below:

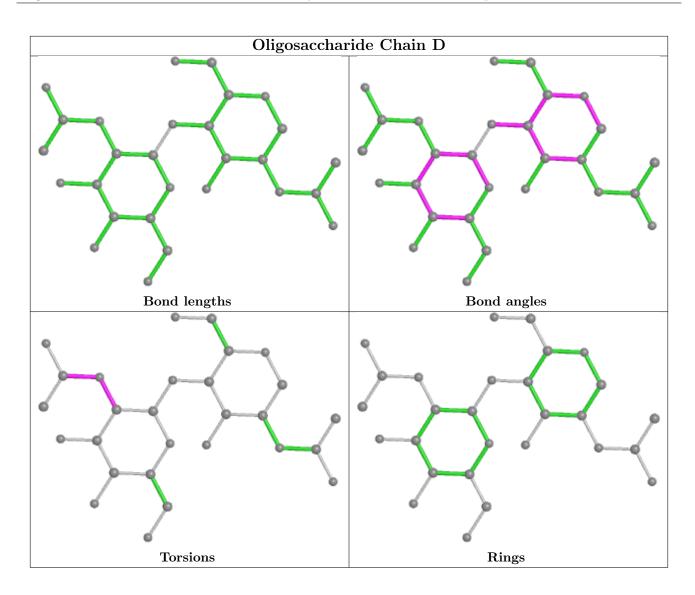
Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C1-C2-N2-C7
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2
4	D	2	NAG	C3-C2-N2-C7

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	Trino	Chain	Res	Link	Bond lengths			В	Bond angles		
Wioi Ty	Type			LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
5	NAG	В	201	2	14,14,15	0.65	0	17,19,21	1.08	1 (5%)	
6	EDO	В	202	-	3,3,3	0.51	0	2,2,2	0.20	0	
5	NAG	A	500	1	14,14,15	0.44	0	17,19,21	0.76	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
5	NAG	В	201	2	-	4/6/23/26	0/1/1/1
6	EDO	В	202	-	-	1/1/1/1	-
5	NAG	A	500	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
5	В	201	NAG	C4-C3-C2	2.19	114.23	111.02

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	В	201	NAG	C8-C7-N2-C2
5	В	201	NAG	O7-C7-N2-C2
5	В	201	NAG	O5-C5-C6-O6
5	В	201	NAG	C4-C5-C6-O6
5	A	500	NAG	C8-C7-N2-C2
5	A	500	NAG	O7-C7-N2-C2
6	В	202	EDO	O1-C1-C2-O2

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^{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>	$\# \mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	A	179/189 (94%)	-0.32	1 (0%) 89 88	14, 24, 50, 81	0
2	В	189/200 (94%)	-0.21	5 (2%) 56 54	14, 25, 53, 87	0
3	С	12/13 (92%)	-0.30	0 100 100	19, 26, 45, 62	0
All	All	380/402 (94%)	-0.27	6 (1%) 72 70	14, 25, 53, 87	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	108	PRO	4.6
2	В	107	GLN	3.4
2	В	110	GLN	3.2
1	A	51	PHE	3.0
2	В	111	HIS	2.9
2	В	109	LEU	2.7

6.2 Non-standard residues in protein, DNA, RNA chains (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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	CIR	С	6	11/12	0.98	0.07	15,16,26,29	0

6.3 Carbohydrates (i)

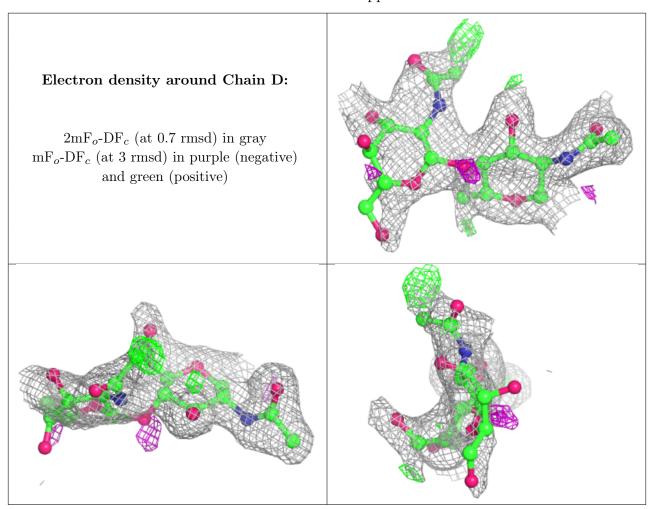
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



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.

M	Iol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	4	NAG	D	2	14/15	0.69	0.31	77,82,86,89	0
4	4	NAG	D	1	14/15	0.88	0.18	24,40,63,75	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^{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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
5	NAG	В	201	14/15	0.77	0.32	64,75,83,84	0



 $Continued\ from\ previous\ page...$

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
5	NAG	A	500	14/15	0.83	0.21	35,57,62,65	0
6	EDO	В	202	4/4	0.90	0.24	47,52,53,55	0

6.5 Other polymers (i)

There are no such residues in this entry.

