

# Full wwPDB X-ray Structure Validation Report (i)

#### Oct 23, 2021 – 12:11 PM EDT

PDB ID : 1FZM

Title : MHC CLASS I NATURAL MUTANT H-2KBM8 HEAVY CHAIN COM-

PLEXED WITH BETA-2 MICROGLOBULIN AND VESICULAR STOM-

ATITIS VIRUS NUCLEOPROTEIN

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Deposited on : 2000-10-03

Resolution : 1.80 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 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 : 2.23.2

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

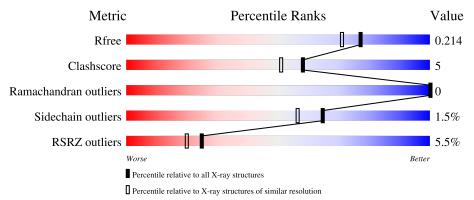
Validation Pipeline (wwPDB-VP) : 2.23.2

## 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 1.80 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	274	6%	89%	11%
2	В	99	5%	89%	11%
3	Р	8		100%	
4	С	3	33%	67%	

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:

$\mathbf{M}$	ol Type	e Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	С	2	-	-	-	X



# 2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 3540 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 H-2 CLASS I HISTOCOMPATIBILITY ANTIGEN, K-B ALPHA CHAIN.

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
1	A	274	Total 2273	C 1432	N 403	O 430	S 8	0	6	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Residue   Modelled   Actual		Comment	Reference
A	22	PHE	TYR	engineered mutation	UNP P01901
A	23	ILE	MET	engineered mutation	UNP P01901
A	24	SER	GLU	engineered mutation	UNP P01901
A	30	ASN ASP		engineered mutation	UNP P01901
A	121	CSO	CYS	modified residue	UNP P01901

• Molecule 2 is a protein called PROTEIN (BETA-2-MICROGLOBULIN).

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	В	99	Total 837	C 533	N 142	O 155	S 7	0	2	0

• Molecule 3 is a protein called PROTEIN (NUCLEOCAPSID PROTEIN).

Mol	Chain	Residues	1	Atoms		ZeroOcc	AltConf	Trace	
3	Р	8	Total 68	C 44	N 12	O 12	0	0	0

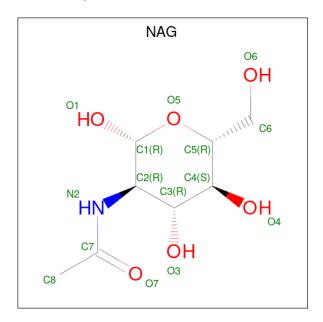
• Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[be ta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.





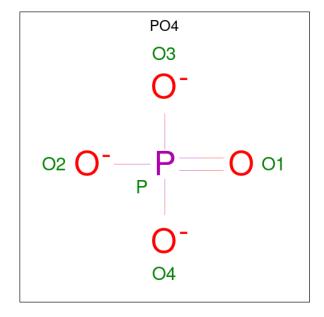
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
4	С	3	Total 38	C 22	N 2	O 14	0	0	0

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



Mo	ol	Chain	Residues	A	ton	ns		ZeroOcc	AltConf
5		Λ	1	Total	С	N	О	0	0
3		Λ	1	14	8	1	5	0	0

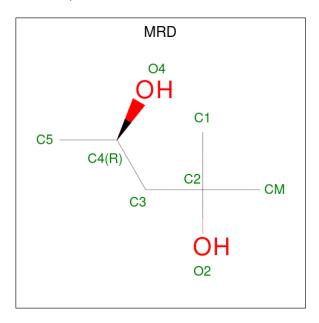
• Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula:  $O_4P$ ).





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

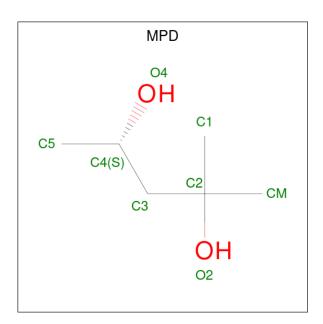
• Molecule 7 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula:  $C_6H_{14}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total 8	C 6	O 2	0	0

• Molecule 8 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:  $C_6H_{14}O_2$ ).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 8 6 2	0	0
8	В	1	Total C O 8 6 2	0	0

### • Molecule 9 is water.

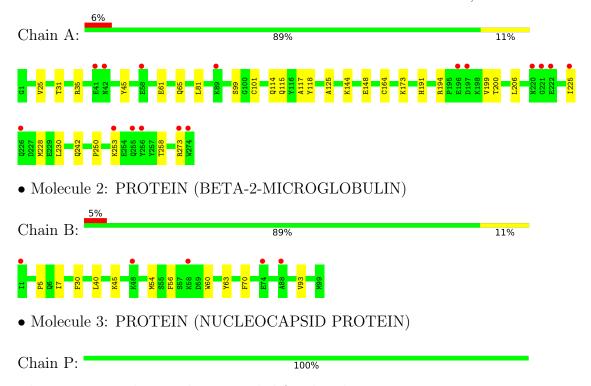
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	196	Total O 196 196	0	0
9	В	73	Total O 73 73	0	0
9	Р	7	Total O 7 7	0	0



## 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: H-2 CLASS I HISTOCOMPATIBILITY ANTIGEN, K-B ALPHA CHAIN



There are no outlier residues recorded for this chain.

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





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	137.26Å 88.15Å 45.59Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.07 - 1.80	Depositor
rtesolution (A)	22.07 - 1.80	EDS
% Data completeness	91.5 (22.07-1.80)	Depositor
(in resolution range)	91.7 (22.07-1.80)	EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.20 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
P. P.	0.210 , $0.223$	Depositor
$R, R_{free}$	0.200 , $0.214$	DCC
$R_{free}$ test set	3769 reflections (7.59%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.0	Xtriage
Anisotropy	0.196	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.41, 56.7	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.48, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3540	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: MPD, CSO, NAG, MRD, PO4, FUL

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		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.35	0/2326	0.62	0/3156	
2	В	0.36	0/864	0.65	0/1171	
3	Р	0.47	0/69	0.78	0/90	
All	All	0.36	0/3259	0.63	0/4417	

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	2273	0	2165	24	0
2	В	837	0	806	5	0
3	Р	68	0	67	0	0
4	С	38	0	34	5	0
5	A	14	0	13	0	0
6	A	5	0	0	0	0
6	В	5	0	0	0	0
7	A	8	0	14	2	0
8	A	8	0	14	0	0
8	В	8	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	196	0	0	2	0
9	В	73	0	0	0	0
9	Р	7	0	0	0	0
All	All	3540	0	3127	31	0

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is 5.

All (31) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

A 4 1	A 4 0	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}\ (\mathring{\rm A})$	overlap (Å)
1:A:173:LYS:HE2	4:C:3:FUL:H62	1.63	0.81
9:A:986:HOH:O	4:C:3:FUL:H2	1.78	0.81
1:A:258:THR:HG22	1:A:273:ARG:HG2	1.65	0.76
1:A:173:LYS:CE	4:C:3:FUL:H62	2.22	0.70
7:A:701:MRD:O2	7:A:701:MRD:H5C3	2.01	0.59
1:A:250:PRO:HG2	1:A:253:LYS:HG2	1.85	0.58
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.39	0.57
1:A:81:LEU:HD23	1:A:118:TYR:CD1	2.41	0.54
1:A:115:GLN:HG2	1:A:125:ALA:HB1	1.89	0.54
1:A:144:LYS:O	1:A:148:GLU:HG3	2.09	0.53
1:A:191:HIS:NE2	1:A:199:VAL:HG11	2.24	0.53
1:A:258:THR:HG22	1:A:273:ARG:CG	2.38	0.53
1:A:31:THR:HG22	9:A:929:HOH:O	2.08	0.53
1:A:225:ILE:HA	1:A:228:MET:CE	2.39	0.52
1:A:25:VAL:HG22	7:A:701:MRD:H5C1	1.92	0.52
1:A:250:PRO:HG2	1:A:253:LYS:CG	2.41	0.51
4:C:1:NAG:O6	4:C:3:FUL:H63	2.10	0.51
1:A:194:ARG:HD2	1:A:200:THR:OG1	2.10	0.51
1:A:61:GLU:O	1:A:65:GLN:HG3	2.13	0.49
1:A:191:HIS:CD2	1:A:199:VAL:HG11	2.49	0.48
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.97	0.46
2:B:7:ILE:HB	2:B:93:VAL:HG21	1.98	0.46
1:A:101:CYS:HG	1:A:164:CYS:CB	2.30	0.44
1:A:99:SER:OG	1:A:114:GLN:NE2	2.51	0.44
1:A:225:ILE:HA	1:A:228:MET:HE2	1.99	0.43
1:A:101:CYS:SG	1:A:164:CYS:SG	3.01	0.43
1:A:225:ILE:HA	1:A:228:MET:HE3	1.99	0.43
2:B:40:LEU:HD23	2:B:45:LYS:HA	2.01	0.43
1:A:173:LYS:HE3	4:C:3:FUL:H62	2.01	0.42
1:A:206:LEU:HD23	1:A:242:GLN:HG3	2.03	0.41

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Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$egin{aligned}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{aligned}$	
2:B:54:MET:HA	2:B:63:TYR:O	2.21	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	Favoured	Allowed	Outliers	Perce	$\mathbf{ntiles}$
1	A	277/274~(101%)	271 (98%)	6 (2%)	0	100	100
2	В	99/99~(100%)	98 (99%)	1 (1%)	0	100	100
3	Р	6/8 (75%)	6 (100%)	0	0	100	100
All	All	382/381 (100%)	375 (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	237/231 (103%)	234 (99%)	3 (1%)	69 62		
2	В	96/94 (102%)	94 (98%)	2 (2%)	53 42		
3	Р	6/6 (100%)	6 (100%)	0	100 100		
All	All	339/331 (102%)	334 (98%)	5 (2%)	65 56		



A 11	/ <b>~</b> \	• 1	• . 1			. 1 1 .		1. / 1	1 1
$A\Pi$	(5)	residiles	with	a	non-rotameric	sidechair	ı are	listed	below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	45	TYR
1	A	230	LEU
2	В	56	PHE
2	В	70	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	114	GLN
1	A	115	GLN
1	A	218	GLN
1	A	220	ASN

#### 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	Chain	Res	Link	B	ond leng	${ m gths}$	В	ond ang	gles
	IVIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2	
Ī	1	CSO	A	121	1	3,6,7	0.55	0	0,6,8	-	-	

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.



$\mathbf{Mol}$	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CSO	A	121	1	-	0/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

### 5.5 Carbohydrates (i)

3 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	Type   Chain   Res   Link		Bo	ond leng	$ ag{ths}$	Bond angles			
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	С	1	1,4	14,14,15	0.57	0	17,19,21	0.70	1 (5%)
4	NAG	С	2	4	14,14,15	0.53	0	17,19,21	0.65	0
4	FUL	С	3	4	10,10,11	0.62	0	14,14,16	0.90	1 (7%)

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	С	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	С	2	4	-	4/6/23/26	0/1/1/1
4	FUL	С	3	4	-	-	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}(^{o})$
4	С	3	FUL	C1-C2-C3	2.11	112.27	109.67
4	С	1	NAG	C2-N2-C7	-2.06	119.96	122.90

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	С	1	NAG	C8-C7-N2-C2
4	С	1	NAG	O7-C7-N2-C2
4	С	2	NAG	C8-C7-N2-C2
4	С	2	NAG	O7-C7-N2-C2
4	С	2	NAG	O5-C5-C6-O6
4	С	2	NAG	C4-C5-C6-O6

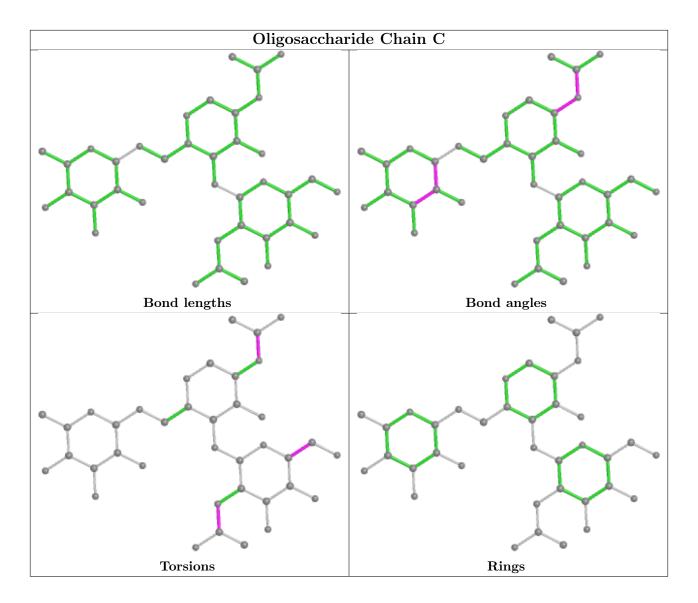
There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	С	1	NAG	1	0
4	С	3	FUL	5	0

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





### 5.6 Ligand geometry (i)

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	Trme	Chain	Res	Link Bond lengths			В	ond ang	cles	
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	PO4	A	801	-	4,4,4	1.65	0	6,6,6	0.43	0
7	MRD	A	701	-	7,7,7	1.26	2 (28%)	9,10,10	0.46	0
6	PO4	В	802	-	4,4,4	1.60	0	6,6,6	0.43	0



Mol	Mol Type Chain Res L		Link	Во	ond leng				cles	
MIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
8	MPD	A	702	-	7,7,7	0.50	0	9,10,10	0.42	0
8	MPD	В	703	-	7,7,7	0.53	0	9,10,10	0.44	0
5	NAG	A	500	1	14,14,15	0.53	0	17,19,21	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	MRD	A	701	-	-	1/5/5/5	-
5	NAG	A	500	1	-	2/6/23/26	0/1/1/1
8	MPD	В	703	-	-	3/5/5/5	-
8	MPD	A	702	-	-	1/5/5/5	-

#### All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	$\operatorname{Ideal}( ext{\AA})$
7	A	701	MRD	C5-C4	-2.25	1.41	1.51
7	A	701	MRD	O4-C4	2.05	1.52	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	500	NAG	C8-C7-N2-C2
5	A	500	NAG	O7-C7-N2-C2
8	В	703	MPD	O2-C2-C3-C4
8	A	702	MPD	O2-C2-C3-C4
7	A	701	MRD	C2-C3-C4-C5
8	В	703	MPD	C1-C2-C3-C4
8	В	703	MPD	CM-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

$\mathbf{Mol}$	Chain	$\operatorname{Res}$	Type	Clashes	Symm-Clashes
7	A	701	MRD	2	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,  $95^{th}$  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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	A	273/274 (99%)	0.12	16 (5%) 22 17	19, 27, 52, 63	0
2	В	99/99 (100%)	0.28	5 (5%) 28 22	21, 29, 40, 49	0
3	Р	8/8 (100%)	0.29	0 100 100	22, 24, 28, 36	0
All	All	380/381 (99%)	0.16	21 (5%) 25 20	19, 27, 49, 63	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	225	ILE	6.1
2	В	1	ILE	5.2
1	A	196	GLU	5.1
1	A	256	TYR	5.1
1	A	220	ASN	5.0
1	A	41	GLU	4.5
1	A	226	GLN	4.1
2	В	48	LYS	3.5
1	A	222	GLU	3.4
1	A	274	TRP	2.9
1	A	253	LYS	2.7
1	A	89	LYS	2.7
2	В	74	GLU	2.7
2	В	88	ALA	2.6
1	A	273	ARG	2.6
2	В	58	LYS	2.6
1	A	221	GLY	2.5
1	A	197	ASP	2.4
1	A	42	ASN	2.2
1	A	58	GLU	2.1
1	A	255	GLN	2.0



#### 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
1	CSO	A	121	7/8	0.92	0.10	21,24,33,35	0

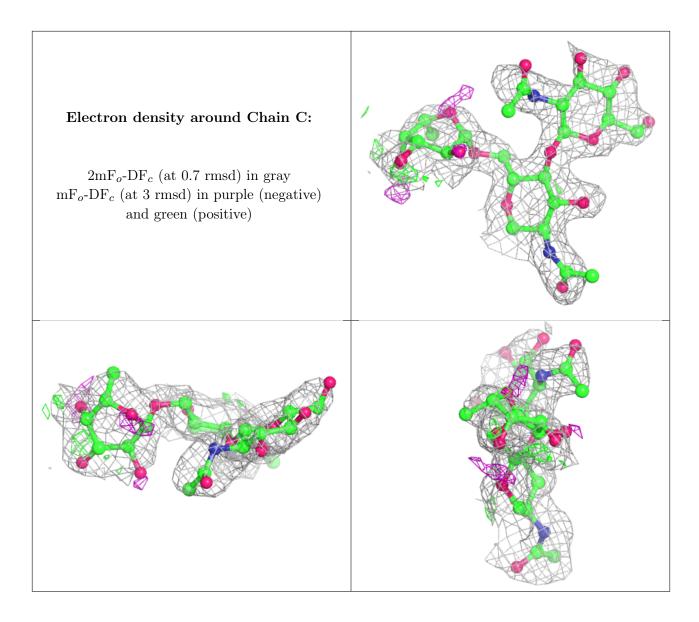
### 6.3 Carbohydrates (i)

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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4	FUL	С	3	10/11	0.67	0.33	56,57,58,58	0
4	NAG	С	2	14/15	0.79	0.43	58,60,61,62	0
4	NAG	С	1	14/15	0.83	0.28	46,51,55,55	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	A	500	14/15	0.76	0.39	50,53,56,56	0
6	PO4	A	801	5/5	0.79	0.25	80,80,80,80	0
8	MPD	В	703	8/8	0.88	0.18	33,38,41,42	0
7	MRD	A	701	8/8	0.91	0.14	30,36,36,37	0
6	PO4	В	802	5/5	0.91	0.15	57,58,59,59	0
8	MPD	A	702	8/8	0.94	0.15	33,36,37,37	0



# 6.5 Other polymers (i)

There are no such residues in this entry.

