

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

#### Mar 23, 2024 – 01:13 PM EDT

PDB ID : 3N41

Title: Crystal structure of the mature envelope glycoprotein complex (spontaneous

cleavage) of Chikungunya virus.

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Deposited on : 2010-05-21

Resolution : 3.01 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 (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:

 $Mol Probity \quad : \quad 4.02b\text{--}467$ 

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36.1

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)

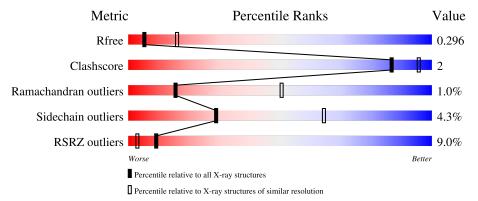
Validation Pipeline (wwPDB-VP) : 2.36.1

# 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 3.01 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\AA)}) \end{array}$
$R_{free}$	130704	2399 (3.04-3.00)
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)
RSRZ outliers	127900	2287 (3.04-3.00)

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	65	69%	12% •	15%					
2	В	360	85%		9% 6%					
3	F	473	13%	6% •	19%					
4	С	2	100%							



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6070 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 E3 envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	55	Total 427	C 268	N 67	O 83	S 9	0	0	0

• Molecule 2 is a protein called E2 envelope glycoprotein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	В	338	Total 2669	C 1666	N 483	O 500	S 20	0	1	0

• Molecule 3 is a protein called E1 envelope glycoprotein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
3	F	382	Total 2909	C 1838	N 487	O 560	S 24	0	1	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLY	-	expression tag	UNP Q1H8W5

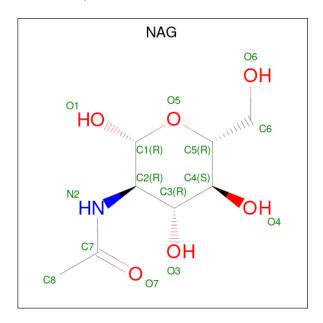
• Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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



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



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

• Molecule 6 is water.

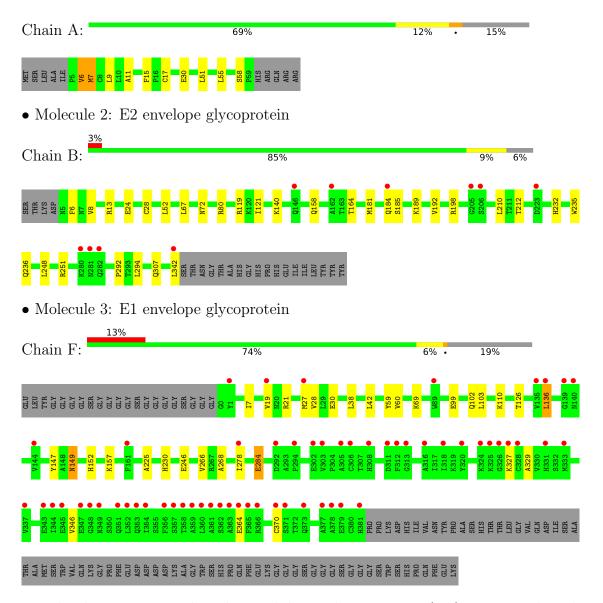
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	В	7	Total O 7 7	0	0
6	F	16	Total O 16 16	0	0



# 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: E3 envelope glycoprotein



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



Chain C: 100%





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	50.24Å 99.61Å 107.10Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $102.90^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	50.00 - 3.01	Depositor
resolution (A)	72.07 - 3.01	EDS
% Data completeness	98.3 (50.00-3.01)	Depositor
(in resolution range)	98.3 (72.07-3.01)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$< I/\sigma(I) > 1$	1.82  (at  3.01Å)	Xtriage
Refinement program	BUSTER 2.9.3	Depositor
$R, R_{free}$	0.252 , $0.275$	Depositor
it, it free	0.272 , $0.296$	DCC
$R_{free}$ test set	1029 reflections $(5.08\%)$	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	70.7	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	$0.32\;,52.7$	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	6070	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.76% 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: NDG, 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		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.45	0/438	0.62	0/598	
2	В	0.37	0/2741	0.60	0/3731	
3	F	0.35	0/2981	0.56	0/4064	
All	All	0.37	0/6160	0.58	0/8393	

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	427	0	410	6	0
2	В	2669	0	2587	8	0
3	F	2909	0	2818	12	0
4	С	28	0	24	0	0
5	F	14	0	13	0	0
6	В	7	0	0	0	0
6	F	16	0	0	0	0
All	All	6070	0	5852	25	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.



The worst 5 of 25 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{ll}  ext{Interatomic} \  ext{distance} & ( ext{Å}) \end{array}$	$egin{array}{c} \operatorname{Clash} \\ \operatorname{overlap}\ ( ext{Å}) \end{array}$	
1:A:6:VAL:HG12	1:A:6:VAL:O	1.71	0.90	
1:A:6:VAL:O	1:A:6:VAL:CG1	2.30	0.79	
2:B:198:ARG:HA	2:B:210:LEU:H	1.51	0.73	
3:F:225:ALA:H	3:F:230:HIS:HE1	1.38	0.72	
3:F:42:LEU:HD11	3:F:266:VAL:HG22	1.85	0.58	

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	53/65 (82%)	50 (94%)	1 (2%)	2 (4%)	3 17		
2	В	336/360 (93%)	313 (93%)	19 (6%)	4 (1%)	13 46		
3	F	381/473 (80%)	369 (97%)	10 (3%)	2 (0%)	29 66		
All	All	770/898 (86%)	732 (95%)	30 (4%)	8 (1%)	15 50		

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	В	121	ILE
3	F	99	GLU
1	A	11	ALA
2	В	72	ASN
2	В	164	THR

#### 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	52/61 (85%)	47 (90%)	5 (10%)	8 30		
2	В	300/319~(94%)	287 (96%)	13 (4%)	29 64		
3	F	319/378 (84%)	308 (97%)	11 (3%)	37 72		
All	All	671/758 (88%)	642 (96%)	29 (4%)	29 64		

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	В	235	TRP
3	F	364	GLU
2	В	342	LEU
3	F	284	GLU
2	В	307	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type	
3	F	149	ASN	
3	F	152	HIS	
3	F	230	HIS	
2	В	236	GLN	
2	В	5	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)

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

## 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	Trme	Chain	Dag	Link	Bo	Bond lengths			Bond angles		
MIOI	Type		Res		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
4	NAG	С	1	4,2	14,14,15	1.38	1 (7%)	17,19,21	1.56	3 (17%)	
4	NDG	С	2	4	14,14,15	1.27	1 (7%)	17,19,21	1.72	6 (35%)	

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	4,2	-	0/6/23/26	0/1/1/1
4	NDG	С	2	4	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
4	С	1	NAG	C1-C2	3.46	1.57	1.52
4	С	2	NDG	C1-C2	2.66	1.56	1.52

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
4	С	1	NAG	C1-O5-C5	3.39	116.78	112.19
4	С	2	NDG	C1-O5-C5	3.25	116.60	112.19
4	С	1	NAG	O4-C4-C3	3.14	117.61	110.35
4	С	2	NDG	O5-C5-C4	3.08	118.32	110.83
4	С	2	NDG	C3-C4-C5	2.60	114.87	110.24

There are no chirality outliers.

All (2) torsion outliers are listed below:

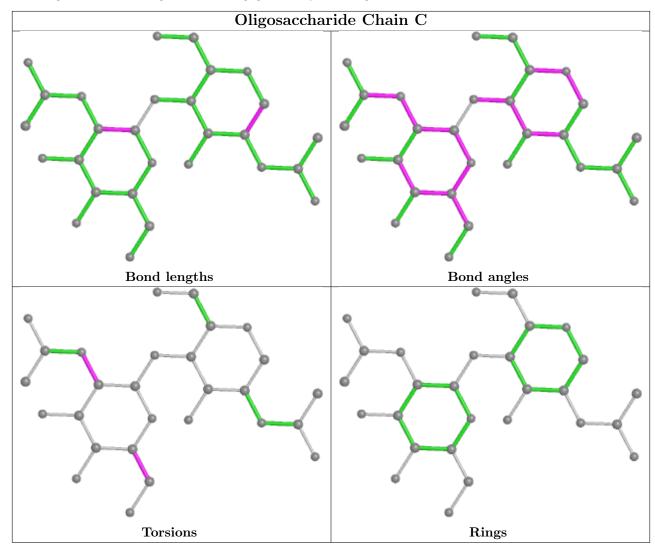
Mol	Chain	Res	Type	Atoms
4	С	2	NDG	C3-C2-N2-C7
4	С	2	NDG	C4-C5-C6-O6



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.



### 5.6 Ligand geometry (i)

#### 1 ligand 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	Pos	Link	Bo	ond leng	ths	В	ond ang	les
IVIOI			nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
5	NAG	F	1001	3	14,14,15	1.30	2 (14%)	17,19,21	1.19	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	F	1001	3	-	2/6/23/26	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\mathring{\mathbf{A}})$	Ideal(Å)
5	F	1001	NAG	C1-C2	3.05	1.56	1.52
5	F	1001	NAG	C3-C2	2.20	1.57	1.52

#### All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
5	F	1001	NAG	C4-C3-C2	3.18	115.67	111.02

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	F	1001	NAG	C1-C2-N2-C7
5	F	1001	NAG	C4-C5-C6-O6

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 > 2	$OWAB(Å^2)$	Q < 0.9
1	A	55/65~(84%)	0.13	0 100 100	63, 81, 92, 101	0
2	В	338/360 (93%)	0.25	10 (2%) 50 22	45, 77, 109, 155	0
3	F	382/473 (80%)	0.79	60 (15%) 2 0	48, 85, 165, 177	0
All	All	775/898 (86%)	0.51	70 (9%) 9 3	45, 81, 163, 177	0

The worst 5 of 70 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	326	GLY	10.7
3	F	327	LYS	7.2
3	F	379	GLU	6.3
3	F	365	PHE	6.1
3	F	360	LEU	6.0

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

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

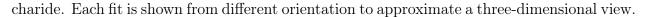
### 6.3 Carbohydrates (i)

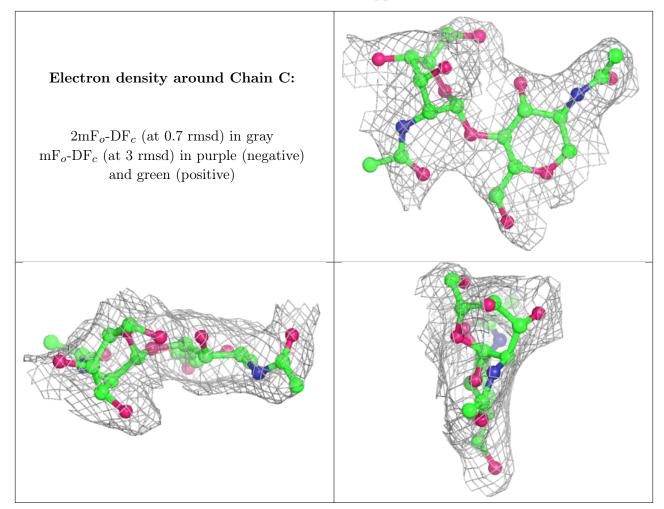
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median,  $95^{th}$  percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

M	ol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4		NDG	С	2	14/15	0.82	0.27	85,89,91,91	0
4		NAG	С	1	14/15	0.88	0.16	78,82,85,85	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-







## 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	${f B-factors}({f \AA}^2)$	Q<0.9
5	NAG	F	1001	14/15	0.60	0.35	115,119,121,121	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

