

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

#### Oct 30, 2023 - 09:05 PM JST

PDB ID	:	4ZFY
Title	:	Structural studies on a non-toxic homologue of type II RIPs from Momordica
		charantia (bitter gourd) in complex with ALPHA-METHYL-D-GALACTOSI
		DE
Authors	:	Chandran, T.; Sharma, A.; Vijayan, M.
Deposited on		
Resolution	:	2.42  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:

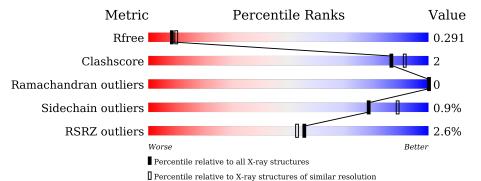
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as 541 be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal 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:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 2.42 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	4647(2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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	А	247	% 94%	6%
2	В	261	4% 97%	•
3	С	3	67% 33%	



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 4154 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 rRNA N-glycosidase.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	А	247	Total 1944	C 1248	N 308	O 377	S 11	0	0	0

• Molecule 2 is a protein called rRNA N-glycosidase.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	В	261	Total 2030	C 1253	N 361	O 402	S 14	0	0	0

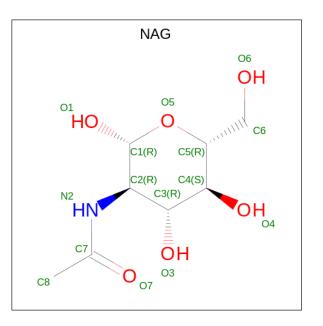
• Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b eta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	I	Aton	ns		ZeroOcc	AltConf	Trace
3	С	3	Total 39	C 22	N 2	O 15	0	0	0

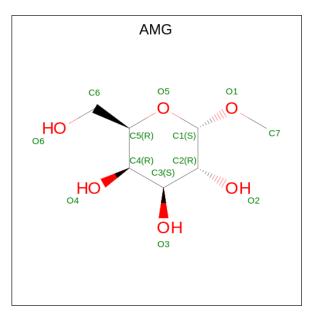
• Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	А	1	Total	С	N	0	0	0
			14	8	1	б		

- Molecule 5 is methyl alpha-D-galactopyranoside (three-letter code: AMG) (formula:  $\rm C_7H_{14}O_6).$ 



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	В	1	Total 13	С 7	O 6	0	0

• Molecule 6 is water.

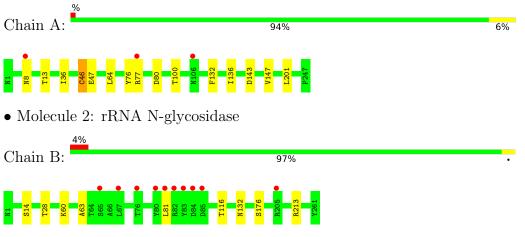


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	44	Total O 44 44	0	0
6	В	70	Total         O           70         70	0	0



# 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: rRNA N-glycosidase

• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-ac etamido-2-deoxy-beta-D-glucopyranose

Chain C:	67%	33%
NAG1 NAG2 BMA3		



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	110.30Å 136.99Å 44.40Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	58.19 - 2.42	Depositor
Resolution (A)	58.19 - 2.42	EDS
% Data completeness	96.3 (58.19-2.42)	Depositor
(in resolution range)	96.3 (58.19-2.42)	EDS
R <sub>merge</sub>	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.44 (at 2.42 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
D D.	0.221 , $0.287$	Depositor
$R, R_{free}$	0.251 , $0.291$	DCC
$R_{free}$ test set	1289 reflections $(5.05\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	36.4	Xtriage
Anisotropy	0.795	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.36 , 23.0	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.45, \langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4154	wwPDB-VP
Average B, all atoms $(Å^2)$	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.15% of the height of the origin peak. No significant pseudotranslation is detected.

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



<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: BMA, NAG, AMG

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	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.45	1/1988~(0.1%)	0.55	0/2701	
2	В	0.36	0/2066	0.60	0/2810	
All	All	0.41	1/4054~(0.0%)	0.58	0/5511	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	47	GLU	CD-OE1	-6.11	1.19	1.25

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	А	1944	0	1888	11	1
2	В	2030	0	1967	4	0
3	С	39	0	34	0	0
4	А	14	0	13	1	0
5	В	13	0	14	0	0
6	А	44	0	0	0	0
6	В	70	0	0	0	0
All	All	4154	0	3916	16	1



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

Atom-1	Atom-2	Interatomic	Clash
		distance $(Å)$	overlap (Å)
2:B:28:THR:O	2:B:60:LYS:HE3	1.70	0.92
1:A:77:ARG:CD	1:A:100:THR:HG21	2.11	0.81
2:B:14:SER:O	2:B:176:SER:HB3	1.91	0.70
1:A:77:ARG:HD3	1:A:100:THR:HG21	1.78	0.65
1:A:8:ASN:O	1:A:13:THR:HG21	1.97	0.65
4:A:301:NAG:O7	4:A:301:NAG:O3	2.17	0.60
1:A:77:ARG:HD2	1:A:100:THR:CG2	2.33	0.59
1:A:80:ASP:O	1:A:100:THR:HB	2.04	0.57
1:A:36:ILE:HD12	1:A:201:LEU:HD21	1.86	0.57
1:A:77:ARG:CD	1:A:100:THR:CG2	2.83	0.55
1:A:77:ARG:HD2	1:A:100:THR:HG21	1.88	0.55
2:B:63:ALA:HB2	2:B:81:LEU:HD11	1.95	0.47
1:A:143:ASP:O	1:A:147:VAL:HG23	2.18	0.43
2:B:28:THR:O	2:B:60:LYS:CE	2.56	0.42
1:A:132:PHE:CE2	1:A:136:ILE:HD11	2.55	0.41
1:A:64:LEU:HD23	1:A:76:TYR:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance} \ (\text{\AA}) \end{array}$	Clash overlap (Å)	
1:A:46:CYS:CB	1:A:46:CYS:SG[2_665]	1.79	0.41	

### 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	А	245/247~(99%)	238~(97%)	7 (3%)	0	100 100	

Continued on next page...



	<i>,</i>	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
2	В	259/261~(99%)	255~(98%)	4 (2%)	0	100	100
All	All	504/508~(99%)	493 (98%)	11 (2%)	0	100	100

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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	А	217/217~(100%)	216 (100%)	1 (0%)	88 95		
2	В	229/229~(100%)	226~(99%)	3~(1%)	69 83		
All	All	446/446 (100%)	442 (99%)	4 (1%)	78 89		

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

Mol	Chain	Res	Type
1	А	46	CYS
2	В	116	THR
2	В	132	ASN
2	В	213	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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)

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 Chain		Chain	Chain	Chain	Chain	Chain	Chain	Chain	Dog	Res	Link	Bo	ond leng	$\mathbf{ths}$	B	ond ang	les
10101	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2							
3	NAG	С	1	3,2	14,14,15	0.52	0	17,19,21	0.78	0							
3	NAG	С	2	3	14,14,15	0.36	0	17,19,21	1.34	4 (23%)							
3	BMA	С	3	3	11,11,12	0.20	0	15,15,17	0.66	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
3	NAG	С	1	$^{3,2}$	-	2/6/23/26	0/1/1/1
3	NAG	С	2	3	-	0/6/23/26	0/1/1/1
3	BMA	$\mathbf{C}$	3	3	_	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	С	2	NAG	C1-O5-C5	3.44	116.85	112.19
3	С	2	NAG	C4-C3-C2	-2.39	107.51	111.02
3	С	2	NAG	C3-C4-C5	-2.15	106.41	110.24
3	С	2	NAG	C2-N2-C7	-2.07	119.95	122.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

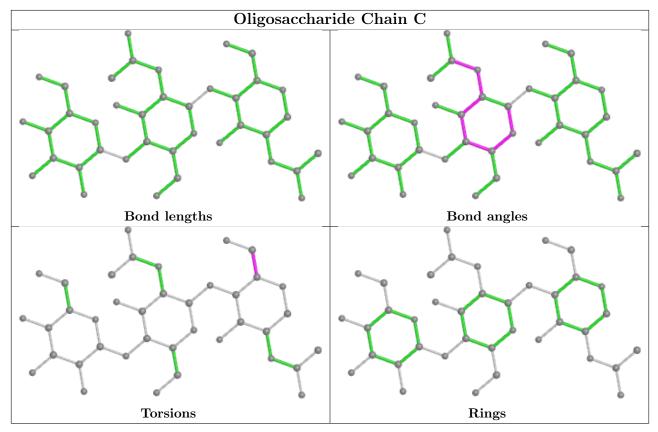
Mol	Chain	Res	Type	Atoms
3	С	1	NAG	O5-C5-C6-O6
3	С	1	NAG	C4-C5-C6-O6



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.



## 5.6 Ligand geometry (i)

2 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).

Mal	Mol Type C	Chain Dec	Link	Bond lengths			Bond angles			
	туре	Chain	nes	S LINK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
4	NAG	А	301	1	14,14,15	0.28	0	17,19,21	0.61	0
5	AMG	В	301	-	13,13,13	0.55	0	18,18,18	0.61	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	А	301	1	-	2/6/23/26	0/1/1/1
5	AMG	В	301	-	-	0/4/24/24	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	А	301	NAG	C1-C2-N2-C7
4	А	301	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	А	301	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,  $95^{th}$  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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	247/247~(100%)	0.22	3 (1%) 79 76	11, 17, 27, 43	0
2	В	261/261~(100%)	0.30	10 (3%) 40 38	11, 16, 27, 44	0
All	All	508/508~(100%)	0.26	13 (2%) 56 53	11, 17, 27, 44	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	82	ARG	3.9
2	В	83	TYR	3.9
1	А	8	ASN	3.6
2	В	85	ASP	3.0
2	В	84	ASP	3.0
2	В	80	TYR	2.6
2	В	81	LEU	2.5
1	А	77	ARG	2.4
2	В	205	ARG	2.3
1	А	106	ASN	2.3
2	В	76	THR	2.1
2	В	67	LEU	2.0
2	В	65	SER	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)

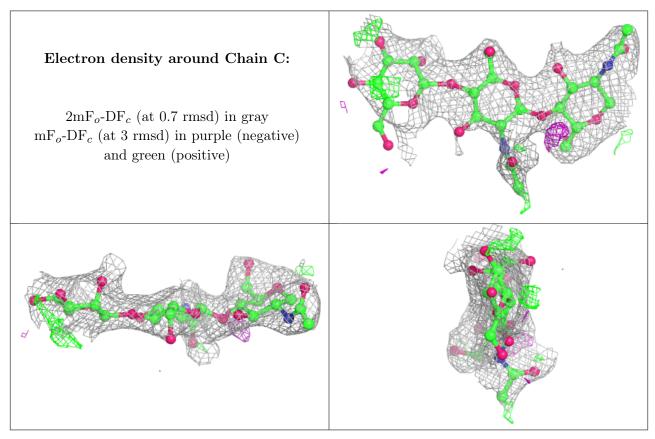
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{B-factors}(\operatorname{\AA}^2)$	Q < 0.9
3	BMA	С	3	11/12	0.69	0.25	$57,\!66,\!73,\!77$	0
3	NAG	С	2	14/15	0.79	0.21	99,107,113,113	0
3	NAG	С	1	14/15	0.87	0.20	41,46,49,50	0

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.

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	B-factors(Å <sup>2</sup> )	Q < 0.9
4	NAG	А	301	14/15	0.59	0.33	37,38,40,40	14
5	AMG	В	301	13/13	0.94	0.23	32,34,37,38	0



## 6.5 Other polymers (i)

There are no such residues in this entry.

