

wwPDB X-ray Structure Validation Summary Report (i)

May 22, 2020 – 11:57 pm BST

PDB ID : 2DDZ

Title: Protein of Unknown Function from Pyrococcus horikoshi

Authors: Rehse, P.H.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative

(RSGI)

Deposited on : 2006-02-07

Resolution : 2.24 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 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 : 2.11

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

Refmac: 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

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

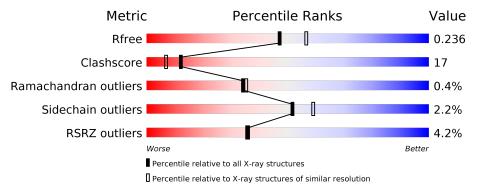
Validation Pipeline (wwPDB-VP) : 2.11

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 2.24 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

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 on 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	190	73%	22%	
1	В	190	71%	27%	
1	С	190	75%	24%	
1	D	190	71%	27%	
1	Е	190	68%	29%	
1	F	190	67%	29%	



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:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GAI	A	514	-	-	X	-
2	GAI	D	513	-	-	X	-
2	GAI	F	509	-	-	X	-



2 Entry composition (i)

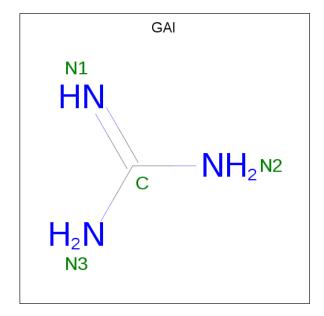
There are 4 unique types of molecules in this entry. The entry contains 9527 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 190aa long hypothetical protein.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	A	188	Total	С	Ν	О	S	0	0	0
1	Λ	100	1507	966	261	275	5	0	0	0
1	В	188	Total	С	N	О	S	0	0	0
1	Ъ	100	1507	966	261	275	5	0	0	0
1	С	188	Total	С	N	О	S	0	0	0
1		100	1507	966	261	275	5	0	0	0
1	D	188	Total	С	N	О	S	0	0	0
1	ש	100	1507	966	261	275	5	0	0	0
1	E	190	Total	С	Ν	О	S	0	0	0
1	ш	190	1523	975	264	278	6	0	U	U
1	F	188	Total	С	N	О	S	0	0	0
1	L'	100	1507	966	261	275	5	U	U	U

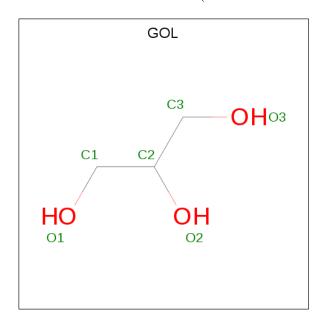
• Molecule 2 is GUANIDINE (three-letter code: GAI) (formula: CH₅N₃).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N 4 1 3	0	0
2	A	1	Total C N 4 1 3	0	0
2	В	1	Total C N 4 1 3	0	0
2	С	1	Total C N 4 1 3	0	0
2	D	1	Total C N 4 1 3	0	0
2	E	1	Total C N 4 1 3	0	0
2	E	1	Total C N 4 1 3	0	0
2	F	1	Total C N 4 1 3	0	0

 \bullet Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $\mathrm{C_3H_8O_3}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	В	1	Total C O 6 3 3	0	0
3	С	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total C O 6 3 3	0	0
3	F	1	Total C O 6 3 3	0	0

• Molecule 4 is water.

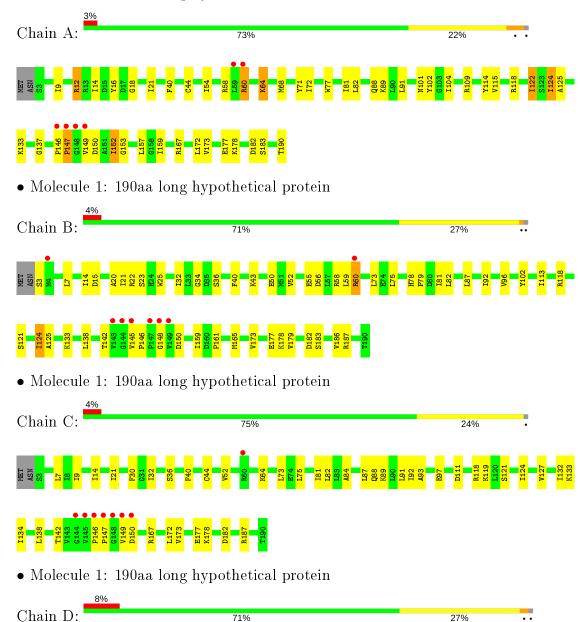
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	83	Total O 83 83	0	0
4	В	65	Total O 65 65	0	0
4	С	76	Total O 76 76	0	0
4	D	47	Total O 47 47	0	0
4	Е	73	Total O 73 73	0	0
4	F	57	Total O 57 57	0	0



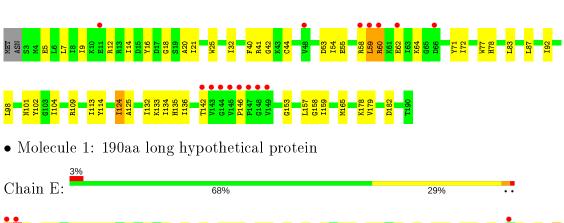
3 Residue-property plots (i)

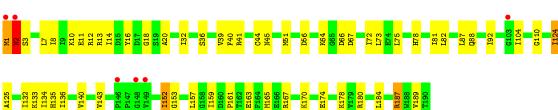
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: 190aa long hypothetical protein

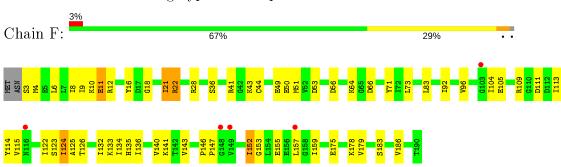








• Molecule 1: 190aa long hypothetical protein





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	$102.76 \text{\AA} 132.43 \text{Å} 135.86 \text{Å}$	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 - 2.24	Depositor
Resolution (A)	48.06 - 2.24	EDS
% Data completeness	92.1 (50.00-2.24)	Depositor
(in resolution range)	92.5 (48.06-2.24)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.56 (at 2.24Å)	Xtriage
Refinement program	CNS	Depositor
D.D.	0.213 , 0.247	Depositor
R, R_{free}	0.204 , 0.236	DCC
R_{free} test set	4480 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.652	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 51.0	EDS
L-test for twinning ²	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9527	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.40% 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: GOL, GAI

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	
WIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5
1	A	0.42	0/1532	0.70	0/2066
1	В	0.45	0/1532	0.70	$1/2066 \ (0.0\%)$
1	С	0.43	0/1532	0.69	0/2066
1	D	0.42	0/1532	0.71	0/2066
1	E	0.44	0/1548	0.71	$2/2087 \; (0.1\%)$
1	F	0.38	0/1532	0.64	0/2066
All	All	0.43	0/9208	0.69	3/12417 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	В	148	GLY	N-CA-C	-6.33	97.26	113.10
1	Е	3	SER	N-CA-C	5.97	127.12	111.00
1	Ε	2	ASN	N-CA-C	5.44	125.69	111.00

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	1507	0	1562	55	0
1	В	1507	0	1562	55	0

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Mol	Chain	Non-H	$\mathbf{H}(\mathbf{model})$	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	С	1507	0	1562	41	0
1	D	1507	0	1562	53	0
1	E	1523	0	1580	73	0
1	F	1507	0	1562	64	0
2	A	8	0	8	2	0
2	В	4	0	4	0	0
2	С	4	0	4	0	0
2	D	4	0	4	2	0
2	E	8	0	8	0	0
2	F	4	0	4	2	0
3	A	6	0	8	0	0
3	В	6	0	8	0	0
3	С	6	0	8	0	0
3	D	6	0	8	1	0
3	E	6	0	8	0	0
3	F	6	0	8	0	0
4	A	83	0	0	0	0
4	В	65	0	0	1	0
4	С	76	0	0	0	0
4	D	47	0	0	0	0
4	Ε	73	0	0	2	0
4	F	57	0	0	0	0
All	All	9527	0	9470	317	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 17.

The worst 5 of 317 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{aligned}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:B:34:GLY:HA3	1:E:1:MET:CG	1.94	0.98
1:C:127:VAL:HG12	1:C:132:ILE:HD13	1.48	0.95
1:D:124:ILE:HD13	1:D:125:ALA:N	1.85	0.92
1:B:124:ILE:HD13	1:B:125:ALA:H	1.32	0.92
1:F:124:ILE:HD13	1:F:125:ALA:N	1.86	0.90

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	Percentiles	
1	A	186/190 (98%)	178 (96%)	7 (4%)	1 (0%)	29	28	
1	В	186/190 (98%)	179 (96%)	6 (3%)	1 (0%)	29	28	
1	С	186/190 (98%)	179 (96%)	7 (4%)	0	100	100	
1	D	186/190 (98%)	177 (95%)	8 (4%)	1 (0%)	29	28	
1	Ε	188/190 (99%)	181 (96%)	6 (3%)	1 (0%)	29	28	
1	F	186/190 (98%)	178 (96%)	8 (4%)	0	100	100	
All	All	1118/1140 (98%)	1072 (96%)	42 (4%)	4 (0%)	34	35	

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	60	ARG
1	E	2	ASN
1	В	60	ARG
1	A	60	ARG

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	Rotameric Outliers	
1	A	166/168~(99%)	158 (95%)	8 (5%)	25 25
1	В	$166/168 \; (99\%)$	165 (99%)	1 (1%)	86 90
1	С	166/168~(99%)	165 (99%)	1 (1%)	86 90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	D	166/168 (99%)	164 (99%)	2 (1%)	71 78		
1	Е	168/168 (100%)	163 (97%)	5 (3%)	41 47		
1	F	166/168 (99%)	161 (97%)	5 (3%)	41 47		
All	All	998/1008 (99%)	976 (98%)	22 (2%)	52 59		

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	167	ARG
1	E	1	MET
1	F	124	ILE
1	D	59	LEU
1	D	124	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	С	116	ASN
1	D	101	ASN
1	F	70	HIS
1	F	116	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)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

14 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	Trans	Chain	Dog	Link	В	Bond lengths			Bond angles		
Mol	Mol Type Chain	$\lim \mid \operatorname{Res} \mid \operatorname{I}$		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2		
2	GAI	В	507	-	3,3,3	0.38	0	3,3,3	0.94	0	
3	GOL	В	502	_	5,5,5	0.34	0	5,5,5	0.30	0	
2	GAI	С	512	_	3,3,3	0.40	0	3,3,3	1.03	0	
2	GAI	Е	511	-	3,3,3	0.32	0	3,3,3	1.02	0	
3	GOL	E	505	_	5,5,5	0.37	0	5,5,5	0.39	0	
3	GOL	С	503	_	5,5,5	0.35	0	5,5,5	0.44	0	
2	GAI	Е	508	_	3,3,3	0.40	0	3,3,3	1.10	0	
2	GAI	A	514	_	3,3,3	0.40	0	3,3,3	1.00	0	
3	GOL	D	504	_	5,5,5	0.39	0	5,5,5	0.46	0	
2	GAI	A	510	_	3,3,3	0.46	0	3,3,3	1.02	0	
2	GAI	D	513	-	3,3,3	0.41	0	3,3,3	1.04	0	
3	GOL	F	506	_	5,5,5	0.40	0	5,5,5	0.34	0	
2	GAI	F	509	-	3,3,3	0.56	0	3,3,3	1.00	0	
3	GOL	A	501	_	5,5,5	0.34	0	5,5,5	0.41	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	\mathbf{Type}	Chain	\mathbf{Res}	Link	Chirals	Torsions	Rings
3	GOL	F	506	_	-	0/4/4/4	-
3	GOL	Ε	505	_	-	0/4/4/4	_
3	GOL	С	503	_	-	0/4/4/4	_
3	GOL	D	504	_	-	0/4/4/4	-
3	GOL	В	502	_	-	0/4/4/4	_
3	GOL	A	501	_	-	0/4/4/4	_

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.



4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	514	GAI	2	0
3	D	504	GOL	1	0
2	D	513	GAI	2	0
2	F	509	GAI	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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	188/190 (98%)	-0.23	6 (3%) 47 47	22, 39, 81, 113	0
1	В	188/190 (98%)	-0.02	8 (4%) 35 34	26, 40, 87, 104	0
1	С	188/190 (98%)	0.01	8 (4%) 35 34	24, 42, 74, 99	0
1	D	188/190 (98%)	0.15	15 (7%) 12 12	28, 52, 94, 118	0
1	E	190/190 (100%)	-0.08	6 (3%) 47 47	24, 43, 73, 108	0
1	F	188/190 (98%)	0.05	5 (2%) 54 55	26, 48, 76, 92	0
All	All	1130/1140 (99%)	-0.02	48 (4%) 36 35	22, 44, 81, 118	0

The worst 5 of 48 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	148	GLY	7.6
1	D	60	ARG	4.8
1	E	1	MET	4.8
1	D	149	VAL	4.6
1	A	148	GLY	4.5

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)

There are no carbohydrates in this entry.



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
2	GAI	E	511	4/4	0.80	0.15	86,91,92,92	0
2	GAI	Е	508	4/4	0.84	0.22	65,69,71,71	0
2	GAI	F	509	4/4	0.84	0.31	72,82,84,86	0
3	GOL	D	504	6/6	0.85	0.15	61,64,68,70	0
2	GAI	A	510	4/4	0.86	0.36	67,71,73,75	0
2	GAI	В	507	4/4	0.87	0.18	62,66,66,68	0
2	GAI	С	512	4/4	0.88	0.17	71,71,75,77	0
2	GAI	D	513	4/4	0.89	0.24	78,87,88,88	0
2	GAI	A	514	4/4	0.89	0.18	71,75,76,76	0
3	GOL	F	506	6/6	0.93	0.15	43,55,64,67	0
3	GOL	В	502	6/6	0.94	0.15	55,60,62,68	0
3	GOL	E	505	6/6	0.95	0.15	44,49,51,54	0
3	GOL	С	503	6/6	0.96	0.13	38,41,46,58	0
3	GOL	A	501	6/6	0.97	0.14	42,55,59,63	0

6.5 Other polymers (i)

There are no such residues in this entry.

