

wwPDB X-ray Structure Validation Summary Report (i)

May 30, 2023 – 01:13 pm BST

PDB ID : 8BCT

Title : X-ray crystal structure of a de novo selected helix-loop-helix heterodimer in a

syn arrangement, 26alpha/26beta

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Deposited on : 2022-10-17

Resolution : 1.70 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 : 4.02b-467

Mogul : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.33

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)

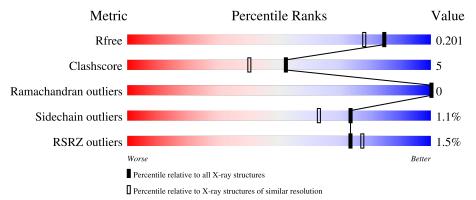
Validation Pipeline (wwPDB-VP) : 2.33

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 1.70 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	В	51	92%	8%	_
1	D	51	88%	10%	
1	F	51	90%	6% •	-
1	Н	51	76%	20% •	-
2	A	51	86%	10% •	•

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Mol	Chain	Length	Quality of chain	
2	С	51	92%	8%
2	Е	51	88%	8% •
2	G	51	88%	6% 6%



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3252 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 26alpha.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	D	50	Total	С	N	О	0	1	0
1	D	30	367	237	61	69	0	1	U
1	В	51	Total	С	N	О	0	1	1
1	Б	91	366	235	61	70	0	1	1
1	Н	49	Total	С	N	O	0	2	0
1	11	49	365	237	61	67	0	2	U
1	F	49	Total	С	N	O	0	3	0
1	1'	49	364	237	60	67	U	ົວ	U

• Molecule 2 is a protein called 26beta.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	G	48	Total	С	N	О	S	0	1	0
	G	40	366	235	59	71	1	0	1	U
2	Е	49	Total	С	N	О	S	0	4	0
	l Li	49	372	240	58	73	1			U
2	Λ	50	Total	С	N	О	S	0	3	0
	Λ	30	386	248	61	76	1		3	U
2	С	51	Total	С	N	О	S	0	2	1
		91	377	242	60	74	1	0	<u> </u>	1

• Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

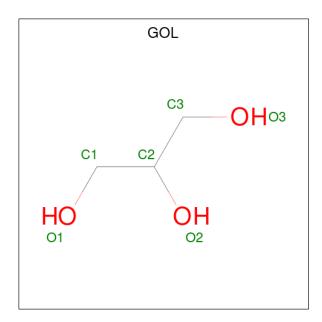




Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total O S 5 4 1	0	0
3	G	1	Total O S 5 4 1	0	0
3	В	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	С	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0

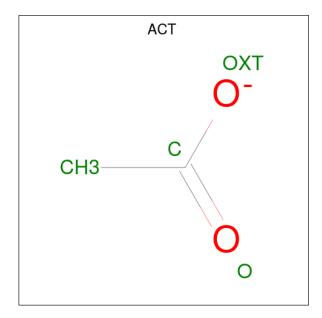
 \bullet Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $\mathrm{C_3H_8O_3}).$





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

 \bullet Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $\mathrm{C_2H_3O_2}).$





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	G	1	Total C O 4 2 2	0	0
5	E	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0

• Molecule 6 is water.

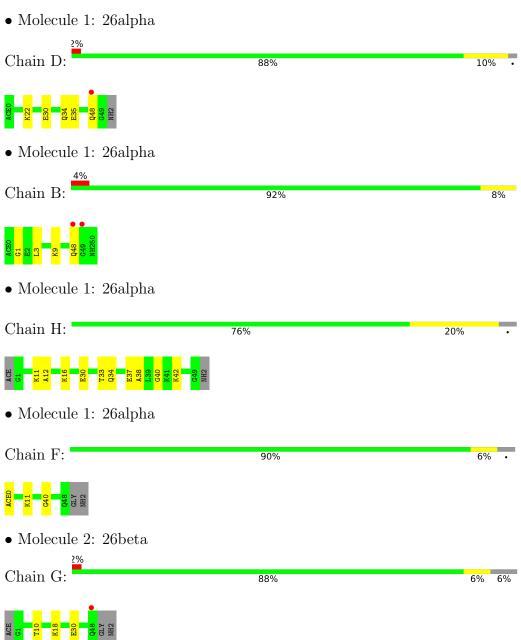
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	32	Total O 32 32	0	0
6	G	23	Total O 23 23	0	0
6	В	30	Total O 30 30	0	0
6	E	29	Total O 29 29	0	0
6	Н	26	Total O 26 26	0	0
6	A	26	Total O 26 26	0	0
6	С	29	Total O 29 29	0	0
6	F	24	Total O 24 24	0	0



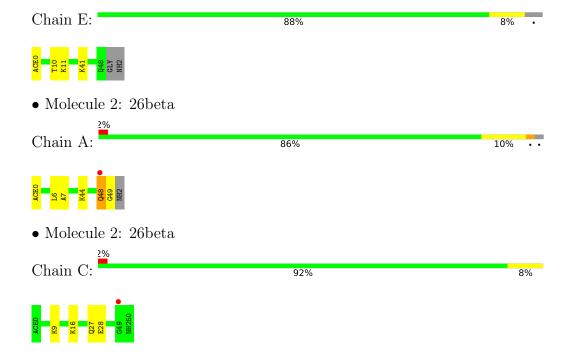
• Molecule 2: 26beta

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.









4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	86.89Å 78.32Å 81.55Å	Depositor
a, b, c, α , β , γ	90.00° 107.97° 90.00°	Depositor
Resolution (Å)	39.16 - 1.70	Depositor
Resolution (A)	39.16 - 1.70	EDS
% Data completeness	99.9 (39.16-1.70)	Depositor
(in resolution range)	99.9 (39.16-1.70)	EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.80 \; (at \; 1.70 \text{Å})$	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.193 , 0.212	Depositor
it, it free	0.203 , 0.201	DCC
R_{free} test set	2808 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.39, 49.8	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3252	wwPDB-VP
Average B, all atoms $(Å^2)$	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 53.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.3187e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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: ACT, ACE, NH2, GOL, SO4

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	\mathbf{angles}
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	В	0.48	0/369	0.69	0/494
1	D	0.56	0/376	0.70	0/501
1	F	0.54	0/379	0.72	0/506
1	Н	0.58	0/379	0.72	0/505
2	A	0.54	0/393	0.76	0/527
2	С	0.65	0/383	0.76	0/513
2	Е	0.68	0/395	0.79	0/529
2	G	0.59	0/372	0.73	0/497
All	All	0.58	0/3046	0.74	0/4072

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	В	366	0	381	8	0
1	D	367	0	395	4	0
1	F	364	0	393	7	0
1	Н	365	0	393	8	0
2	A	386	0	400	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	С	377	0	394	3	0
2	Ε	372	0	391	5	0
2	G	366	0	388	5	0
3	В	5	0	0	0	0
3	С	5	0	0	0	0
3	D	5	0	0	0	0
3	Ε	5	0	0	0	0
3	F	5	0	0	0	0
3	G	5	0	0	0	0
4	В	6	0	8	3	0
4	С	6	0	8	0	0
4	D	6	0	8	3	0
4	Ε	6	0	8	0	0
5	A	4	0	3	0	0
5	Е	4	0	3	0	0
5	F	4	0	3	0	0
5	G	4	0	3	0	0
6	A	26	0	0	0	0
6	В	30	0	0	0	0
6	С	29	0	0	1	0
6	D	32	0	0	0	0
6	Е	29	0	0	0	0
6	F	24	0	0	0	0
6	G	23	0	0	1	0
6	Н	26	0	0	0	0
All	All	3252	0	3179	34	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 5.

The worst 5 of 34 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:B:48:GLN:HB3	1:F:0:ACE:CH3	2.12	0.79
1:B:48:GLN:HB3	1:F:0:ACE:H1	1.66	0.78
1:B:3:LEU:H	4:B:102:GOL:H11	1.48	0.77
2:G:10:THR:HG21	1:H:40:GLY:HA2	1.69	0.75
2:G:18:LYS:HE3	1:H:37:GLU:OE2	1.86	0.74

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	Percei	ntiles
1	В	50/51~(98%)	50 (100%)	0	0	100	100
1	D	$49/51 \ (96\%)$	49 (100%)	0	0	100	100
1	F	50/51~(98%)	47 (94%)	3 (6%)	0	100	100
1	Н	49/51~(96%)	49 (100%)	0	0	100	100
2	A	51/51 (100%)	51 (100%)	0	0	100	100
2	C	51/51~(100%)	51 (100%)	0	0	100	100
2	E	51/51 (100%)	51 (100%)	0	0	100	100
2	G	47/51 (92%)	47 (100%)	0	0	100	100
All	All	398/408 (98%)	395 (99%)	3 (1%)	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	В	33/35~(94%)	33 (100%)	0	100	100	
1	D	35/35~(100%)	34 (97%)	1 (3%)	42	23	
1	F	35/35~(100%)	35 (100%)	0	100	100	
1	Н	34/35~(97%)	34 (100%)	0	100	100	
2	A	$37/37\ (100\%)$	36 (97%)	1 (3%)	44	26	
2	С	$36/37\ (97\%)$	36 (100%)	0	100	100	

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Mol	Chain	Analysed	Rotameric	Rotameric Outliers		Percentiles	
2	E	38/37 (103%)	37 (97%)	1 (3%)	46 28		
2	G	$36/37 \ (97\%)$	36 (100%)	0	100 100)	
All	All	284/288 (99%)	281 (99%)	3 (1%)	73 63		

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	48	GLN
2	Е	11	LYS
2	A	48	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	27	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)

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

5.5 Carbohydrates (i)

There are no monosaccharides 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).

Mol	Trino	Chain	Res	Link	В	ond leng	$_{ m gths}$	Bond angles			
IVIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
3	SO4	С	101	-	4,4,4	0.46	0	6,6,6	0.13	0	
3	SO4	G	101	-	4,4,4	0.46	0	6,6,6	0.06	0	
3	SO4	D	101	-	4,4,4	0.38	0	6,6,6	0.08	0	
3	SO4	Е	101	-	4,4,4	0.44	0	6,6,6	0.06	0	
4	GOL	D	102	-	5,5,5	0.12	0	5,5,5	0.38	0	
3	SO4	В	101	-	4,4,4	0.51	0	6,6,6	0.13	0	
5	ACT	G	102	-	3,3,3	0.94	0	3,3,3	1.45	1 (33%)	
3	SO4	F	101	-	4,4,4	0.39	0	6,6,6	0.58	0	
4	GOL	В	102	-	5,5,5	0.22	0	5,5,5	0.81	0	
5	ACT	A	101	-	3,3,3	1.01	0	3,3,3	0.89	0	
5	ACT	F	102	-	3,3,3	0.95	0	3,3,3	1.02	0	
4	GOL	С	102	-	5,5,5	0.14	0	5,5,5	0.36	0	
5	ACT	Е	102	-	3,3,3	1.51	1 (33%)	3,3,3	0.77	0	
4	GOL	Е	103	-	5,5,5	0.11	0	5,5,5	0.23	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	GOL	С	102	-	-	0/4/4/4	-
4	GOL	D	102	-	-	2/4/4/4	-
4	GOL	Е	103	-	-	2/4/4/4	-
4	GOL	В	102	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mo	l Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
5	E	102	ACT	СН3-С	2.15	1.58	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
5	G	102	ACT	O-C-CH3	-2.01	114.52	122.33

There are no chirality outliers.

5 of 8 torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
4	D	102	GOL	O1-C1-C2-C3
4	В	102	GOL	O1-C1-C2-C3
4	В	102	GOL	C1-C2-C3-O3
4	В	102	GOL	O2-C2-C3-O3
4	В	102	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	102	GOL	3	0
4	В	102	GOL	3	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	В	49/51 (96%)	-0.21	2 (4%) 37 41	27, 39, 53, 65	0
1	D	49/51 (96%)	-0.24	1 (2%) 65 69	23, 36, 57, 61	0
1	F	48/51 (94%)	-0.28	0 100 100	26, 34, 48, 53	0
1	Н	49/51 (96%)	-0.27	0 100 100	22, 35, 47, 75	0
2	A	49/51 (96%)	-0.31	1 (2%) 65 69	26, 34, 49, 71	0
2	С	49/51 (96%)	-0.24	1 (2%) 65 69	21, 33, 53, 73	0
2	E	48/51 (94%)	-0.20	0 100 100	28, 36, 54, 57	0
2	G	48/51 (94%)	-0.26	1 (2%) 63 67	22, 33, 53, 63	0
All	All	389/408 (95%)	-0.25	6 (1%) 73 77	21, 35, 53, 75	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	49	GLY	4.3
2	С	49	GLY	3.5
2	A	48	GLN	3.4
1	В	48	GLN	2.9
1	D	48	GLN	2.4

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 monosaccharides 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	$\operatorname{B-factors}(\AA^2)$	Q<0.9
4	GOL	С	102	6/6	0.78	0.18	51,67,69,75	0
3	SO4	F	101	5/5	0.80	0.16	37,39,47,67	5
5	ACT	F	102	4/4	0.81	0.13	52,59,61,62	0
4	GOL	В	102	6/6	0.82	0.15	43,51,60,68	0
4	GOL	D	102	6/6	0.84	0.18	52,72,77,81	0
3	SO4	С	101	5/5	0.92	0.12	50,54,60,62	5
4	GOL	Е	103	6/6	0.92	0.08	62,68,73,77	0
5	ACT	Е	102	4/4	0.93	0.11	43,45,52,60	0
3	SO4	G	101	5/5	0.93	0.12	46,48,55,57	5
5	ACT	G	102	4/4	0.95	0.07	34,43,45,56	0
3	SO4	В	101	5/5	0.95	0.12	48,53,64,72	5
5	ACT	A	101	4/4	0.95	0.09	65,69,77,78	0
3	SO4	Е	101	5/5	0.95	0.11	49,51,56,58	5
3	SO4	D	101	5/5	0.96	0.16	66,71,74,76	5

6.5 Other polymers (i)

There are no such residues in this entry.

