

wwPDB X-ray Structure Validation Summary Report (i)

Nov 15, 2023 – 02:18 PM JST

PDB ID	:	6JAR
Title	:	Crystal structure of ABC transporter alpha-glycoside-binding mutant protein
		R49A in complex with maltose
Authors	:	Kanaujia, S.P.; Chandravanshi, M.; Gogoi, P.
Deposited on	:	2019-01-24
Resolution	:	1.63 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:

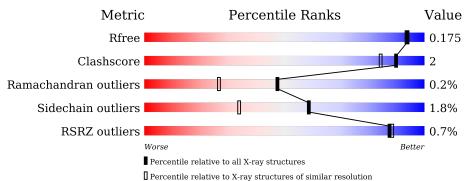
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 1.63 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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	3122(1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079(1.66-1.62)

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	408	% 92%	6% •
2	В	2	100%	



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3910 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 ABC transporter, periplasmic substrate-binding protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	404	Total 3262	C 2073	N 583	O 597	S 9	0	9	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	49	ALA	ARG	engineered mutation	UNP Q5SLD7
А	403	HIS	-	expression tag	UNP Q5SLD7
А	404	HIS	-	expression tag	UNP Q5SLD7
А	405	HIS	-	expression tag	UNP Q5SLD7
А	406	HIS	-	expression tag	UNP Q5SLD7
А	407	HIS	-	expression tag	UNP Q5SLD7
А	408	HIS	-	expression tag	UNP Q5SLD7

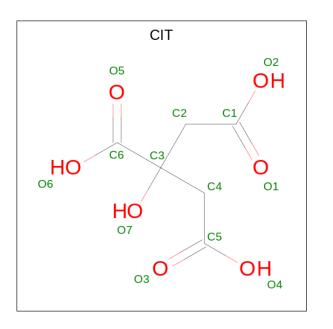
• Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	At	\mathbf{oms}		ZeroOcc	AltConf	Trace
2	В	2	Total 23	C 12	0 11	0	0	0

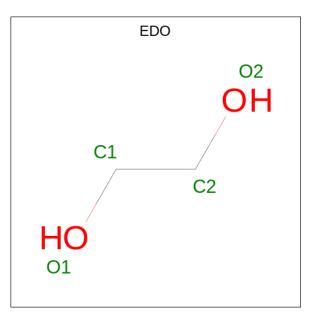
• Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	А	1	Total 13	С 6	O 7	0	0

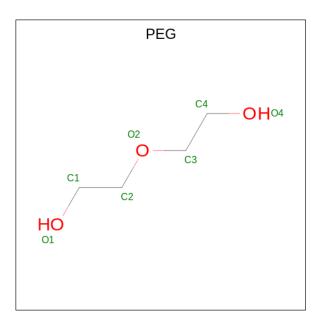
• Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	А	1	Total 4	${ m C} 2$	O 2	0	0

• Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).





Mo	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 7	$\begin{array}{c} \mathrm{C} \\ 4 \end{array}$	O 3	0	0

• Molecule 6 is water.

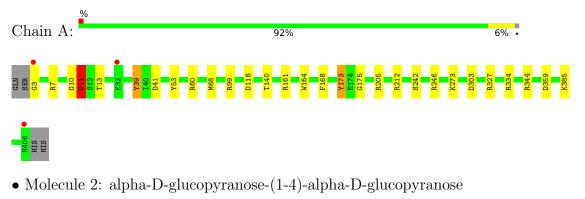
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	601	Total O 601 601	0	0



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: ABC transporter, periplasmic substrate-binding protein



Chain B:

100%

GLC1 GLC2



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	84.98Å 84.98Å 145.46Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.38 - 1.63	Depositor
Resolution (A)	42.11 - 1.63	EDS
% Data completeness	91.1 (73.38-1.63)	Depositor
(in resolution range)	$91.1 \ (42.11 - 1.63)$	EDS
R _{merge}	0.06	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.44 (at 1.63 Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.140 , 0.166	Depositor
II, II, <i>free</i>	0.149 , 0.175	DCC
R_{free} test set	2964 reflections $(4.84%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	16.5	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.33 , 43.4	EDS
L-test for twinning ²	$ \langle L \rangle = 0.49, \langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3910	wwPDB-VP
Average B, all atoms $(Å^2)$	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.62% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: CIT, GLC, PEG, EDO

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	Boi	nd lengths	Bo	ond angles
			# Z > 5	RMSZ	# Z > 5
1	А	1.04	3/3376~(0.1%)	1.08	17/4591~(0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
1	А	11	ASP	C-O	7.36	1.37	1.23
1	А	53	TYR	CE1-CZ	-5.43	1.31	1.38
1	А	60	ARG	CZ-NH2	5.23	1.39	1.33

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	334	ARG	NE-CZ-NH2	-8.29	116.16	120.30
1	А	359	ASP	CB-CG-OD1	7.65	125.19	118.30
1	А	344	ARG	NE-CZ-NH2	-7.53	116.53	120.30
1	А	344	ARG	NE-CZ-NH1	7.46	124.03	120.30
1	А	205	ARG	NE-CZ-NH2	-7.21	116.69	120.30

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	3262	0	3199	12	0
2	В	23	0	21	0	0
3	А	13	0	5	0	0
4	А	4	0	6	0	0
5	А	7	0	10	0	0
6	А	601	0	0	4	2
All	All	3910	0	3241	12	2

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:GLY:O	1:A:39:TYR:OH	1.69	1.11
1:A:303[B]:ASP:OD2	6:A:601:HOH:O	1.89	0.91
1:A:41[B]:ASP:OD2	6:A:602:HOH:O	2.09	0.71
1:A:11:ASP:N	1:A:68[B]:MET:HE3	2.08	0.69
1:A:13:THR:HB	6:A:1059:HOH:O	1.95	0.66

All (2) 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	Interatomic distance (Å)	Clash overlap (Å)
6:A:1086:HOH:O	6:A:1121:HOH:O[7_555]	1.90	0.30
6:A:751:HOH:O	6:A:1025:HOH:O[6_444]	2.17	0.03

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	411/408 (101%)	402 (98%)	8 (2%)	1 (0%)	47 26	



All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	11	ASP

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	337/332~(102%)	331~(98%)	6(2%)	59 34	

5 of 6 residues with a non-rotameric side chain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	173	TYR
1	А	246	ARG
1	А	385	LYS
1	А	118	ASP
1	А	11	ASP

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)

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 Type		e Chain Res Link			Bo	Bond lengths			Bond angles		
	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
2	GLC	В	1	2	12,12,12	1.58	2 (16%)	$17,\!17,\!17$	2.11	<u>6 (35%)</u>	
2	GLC	В	2	2	11,11,12	1.04	0	$15,\!15,\!17$	1.59	4 (26%)	

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.

Mo	ol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2		GLC	В	1	2	-	0/2/22/22	0/1/1/1
2		GLC	В	2	2	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	В	1	GLC	O4-C4	3.79	1.51	1.43
2	В	1	GLC	O3-C3	2.58	1.49	1.43

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	1	GLC	O3-C3-C2	-4.16	100.74	110.35
2	В	1	GLC	O3-C3-C4	3.88	119.31	110.35
2	В	1	GLC	O4-C4-C5	-3.38	100.90	109.30
2	В	1	GLC	C6-C5-C4	-2.97	106.04	113.00
2	В	2	GLC	C1-C2-C3	2.92	113.25	109.67

There are no chirality outliers.

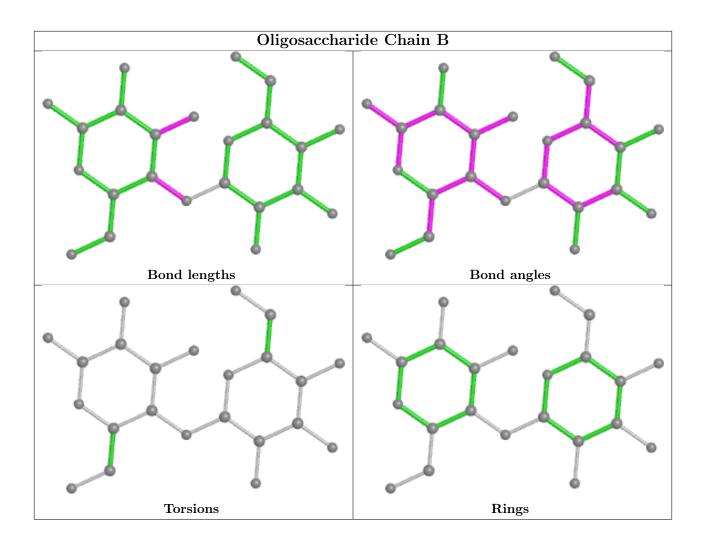
There are no torsion outliers.

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)

3 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	Type	Chain	Res	Link	Bo	Bond lengths			Bond angles		
	01				Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
5	PEG	А	504	-	$6,\!6,\!6$	0.54	0	$5,\!5,\!5$	0.94	0	
3	CIT	А	502	-	12,12,12	1.74	3 (25%)	17,17,17	1.56	4 (23%)	
4	EDO	А	503	-	3,3,3	0.39	0	2,2,2	0.86	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
5	PEG	А	504	-	-	3/4/4/4	-
3	CIT	А	502	-	-	4/16/16/16	-
4	EDO	А	503	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
3	А	502	CIT	C2-C3	-3.87	1.49	1.53
3	А	502	CIT	C4-C3	-2.66	1.50	1.53
3	А	502	CIT	C3-C6	-2.30	1.51	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	502	CIT	O7-C3-C6	2.92	112.96	108.86
3	А	502	CIT	C4-C3-C6	-2.77	104.16	110.11
3	А	502	CIT	O5-C6-C3	-2.64	118.51	122.25
3	А	502	CIT	O3-C5-C4	-2.35	116.06	122.94

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	А	504	PEG	C4-C3-O2-C2
5	А	504	PEG	O1-C1-C2-O2
5	А	504	PEG	O2-C3-C4-O4
3	А	502	CIT	C2-C3-C6-O5
3	А	502	CIT	C2-C3-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	$\langle RSRZ \rangle$	#RSRZ>2	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	А	404/408~(99%)	-0.32	3 (0%) 87 88	10, 16, 31, 64	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	406	HIS	3.7
1	А	3	GLY	2.9
1	А	32	LYS	2.2

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

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

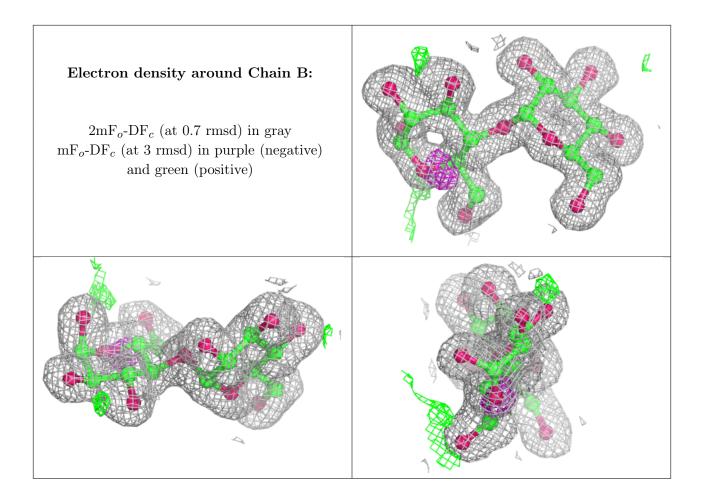
6.3 Carbohydrates (i)

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.

ſ	Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
	2	GLC	В	1	12/12	0.91	0.11	13,23,27,39	0
	2	GLC	В	2	11/12	0.98	0.09	10,11,11,12	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	B-factors(Å ²)	Q<0.9
4	EDO	А	503	4/4	0.72	0.13	41,46,49,54	0
5	PEG	А	504	7/7	0.88	0.25	38,45,50,52	0
3	CIT	А	502	13/13	0.94	0.09	16,24,29,35	0

6.5 Other polymers (i)

There are no such residues in this entry.

