

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

#### Nov 21, 2023 – 07:12 AM JST

:	7F6Q
:	Crystal structure of metal-citrate-binding mutant (S79A) protein (MctA) of
	ABC transporter in apo state
:	Kanaujia, S.P.; Mandal, S.K.
	2021-06-25
:	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:

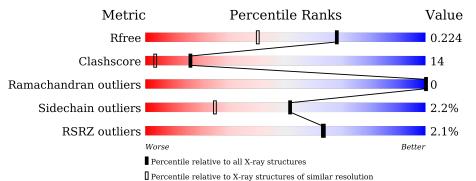
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal 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 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 <sub>free</sub>	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		
			2%		
1	А	342	82%	14% ••	

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
4	CO2	А	406	-	-	Х	-
5	ACT	А	407	-	-	Х	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	EDO	А	409	-	-	Х	-
6	EDO	А	410	-	-	Х	-
6	EDO	А	411	-	-	Х	-
7	PDO	А	412	-	-	Х	-
8	PEG	А	414	-	-	Х	-

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# 2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 3128 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 Iron ABC transporter, periplasmic iron-binding protein.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	338	Total 2726	C 1752	N 472	0 498	$\frac{S}{4}$	0	11	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-1	MET	-	initiating methionine	UNP Q53VZ2
А	0	MET	-	expression tag	UNP Q53VZ2
А	79	ALA	SER	engineered mutation	UNP Q53VZ2

• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

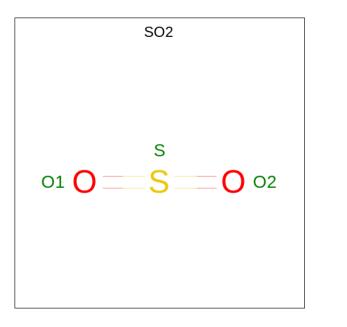
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	4	Total Cl 4 4	0	0

• Molecule 3 is SULFUR DIOXIDE (three-letter code: SO2) (formula: O<sub>2</sub>S).



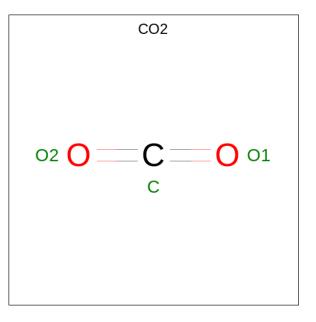
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	А	1	Total 3	0 2	S 1	0	0

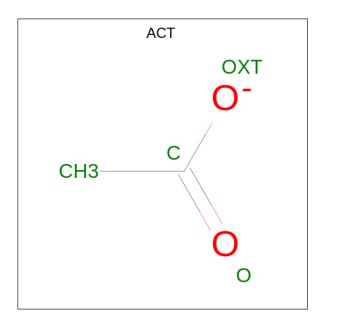
• Molecule 4 is CARBON DIOXIDE (three-letter code: CO2) (formula:  $CO_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	А	1	Total 3	С 1	O 2	0	0

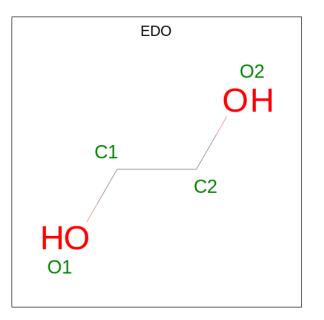
 $\bullet\,$  Molecule 5 is ACETATE ION (three-letter code: ACT) (formula:  $\rm C_2H_3O_2).$ 





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	А	1	Total 4	$\begin{array}{c} \mathrm{C} \\ \mathrm{2} \end{array}$	O 2	0	0

• Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



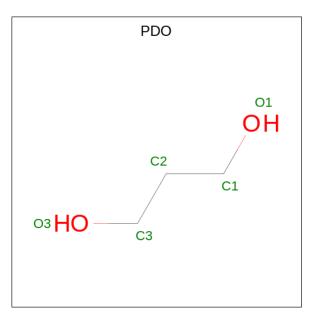
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	A	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0

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Μ	ol	Chain	Residues	Atoms		ZeroOcc	AltConf	
6	<b>j</b>	А	1	Total 4	${ m C} 2$	O 2	0	0

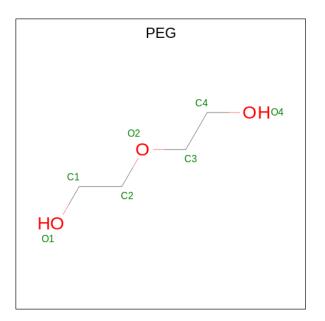
• Molecule 7 is 1,3-PROPANDIOL (three-letter code: PDO) (formula:  $C_3H_8O_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 5 & 3 & 2 \end{array}$	0	0
7	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 5 & 3 & 2 \end{array}$	0	0

• Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
8	А	1	Total 7	$\begin{array}{c} \mathrm{C} \\ 4 \end{array}$	O 3	0	0

• Molecule 9 is water.

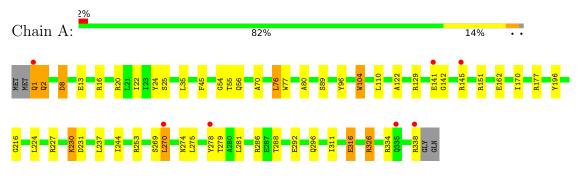
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	А	355	Total C 355 35	5	0	0



# 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: Iron ABC transporter, periplasmic iron-binding protein





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants	72.17Å 72.17Å 114.30Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	42.21 - 1.63	Depositor
Resolution (A)	42.18 - 1.63	EDS
% Data completeness	99.9 (42.21-1.63)	Depositor
(in resolution range)	99.9(42.18-1.63)	EDS
R <sub>merge</sub>	0.06	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$4.30 (at 1.63 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
P. P.	0.175 , $0.213$	Depositor
$R, R_{free}$	0.186 , $0.224$	DCC
$R_{free}$ test set	2174 reflections $(4.97%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	19.7	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.33 , $42.3$	EDS
L-test for twinning <sup>2</sup>	$< L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3128	wwPDB-VP
Average B, all atoms $(Å^2)$	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.70% 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: CO2, ACT, PEG, PDO, CL, EDO, SO2

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	
	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	1.00	5/2815~(0.2%)	1.19	12/3817~(0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1

Ζ Chain Observed(Å) Mol Res Type Atoms Ideal(Å) 1 А 316 GLU CD-OE1 7.331.331.251 А  $\mathbf{2}$  $\operatorname{GLN}$ C-O 6.661.361.231 А 13 GLU CD-OE2 6.511.321.25THR C-O 1 А 2795.361.33 1.23 GLU 1 А 292CD-OE1 -5.161.201.25

All (5) bond length outliers are listed below:

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	253	ARG	NE-CZ-NH2	-10.61	115.00	120.30
1	А	253	ARG	NE-CZ-NH1	9.61	125.11	120.30
1	А	286	ARG	NE-CZ-NH1	7.91	124.26	120.30
1	А	326	ARG	NE-CZ-NH2	6.99	123.79	120.30
1	А	129	ARG	NE-CZ-NH1	6.90	123.75	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:



$\mathbf{M}$	ol	Chain	Res	Type	Group
1		А	269	SER	Mainchain

#### 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	2726	0	2769	71	0
2	А	4	0	0	1	0
3	А	3	0	0	0	0
4	А	3	0	0	3	0
5	А	4	0	3	3	0
6	А	16	0	23	27	0
7	А	10	0	16	12	0
8	А	7	0	9	4	0
9	А	355	0	0	18	0
All	All	3128	0	2820	81	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 14.

The worst 5 of 81 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:274:ASN:HB2	6:A:409:EDO:C1	1.78	1.11
1:A:122:ALA:CB	1:A:170[B]:ILE:HD11	1.84	1.06
1:A:122:ALA:HB2	1:A:170[B]:ILE:HD11	1.40	0.97
1:A:96:TYR:OH	7:A:412:PDO:H31	1.69	0.93
1:A:274:ASN:HB2	6:A:409:EDO:H12	1.49	0.92

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	А	347/342~(102%)	343~(99%)	4 (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	А	281/273~(103%)	274~(98%)	7(2%)	47 20	

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	104	TRP
1	А	110	LEU
1	А	230	LYS
1	А	151	ARG
1	А	76[B]	LEU

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)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry (i)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 for Mogul analysis.

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	Turne	Chain	Res	Link	B	ond leng	gths	В	ond ang	gles
INIOI	Type	Chain	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
6	EDO	А	408	-	3,3,3	0.16	0	$2,\!2,\!2$	0.14	0
4	CO2	А	406	-	2,2,2	0.19	0	$1,\!1,\!1$	0.81	0
3	SO2	А	405	-	2,2,2	0.52	0	$1,\!1,\!1$	0.15	0
7	PDO	А	413	-	4,4,4	0.15	0	$3,\!3,\!3$	0.43	0
8	PEG	А	414	-	6,6,6	0.55	0	$5,\!5,\!5$	0.17	0
6	EDO	А	409	-	3,3,3	0.35	0	$2,\!2,\!2$	0.59	0
6	EDO	А	410	-	3,3,3	0.62	0	$2,\!2,\!2$	0.27	0
7	PDO	А	412	-	4,4,4	0.35	0	$3,\!3,\!3$	0.95	0
5	ACT	А	407	-	3,3,3	1.33	0	$3,\!3,\!3$	0.31	0
6	EDO	А	411	-	3,3,3	0.85	0	$2,\!2,\!2$	0.67	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
6	EDO	А	408	-	-	1/1/1/1	-
7	PDO	А	413	-	-	0/2/2/2	-
8	PEG	А	414	-	-	2/4/4/4	-
6	EDO	А	409	-	-	1/1/1/1	-
6	EDO	А	410	-	-	1/1/1/1	-
7	PDO	А	412	-	-	2/2/2/2	-
6	EDO	А	411	-	-	1/1/1/1	_

There are no bond length outliers.



There are no bond angle outliers.

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	А	412	PDO	O1-C1-C2-C3
7	А	412	PDO	C1-C2-C3-O3
8	А	414	PEG	O2-C3-C4-O4
6	А	408	EDO	O1-C1-C2-O2
8	А	414	PEG	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	А	408	EDO	2	0
4	А	406	CO2	3	0
7	А	413	PDO	2	0
8	А	414	PEG	4	0
6	А	409	EDO	10	0
6	А	410	EDO	11	0
7	А	412	PDO	10	0
5	А	407	ACT	3	0
6	А	411	EDO	4	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>2		$OWAB(Å^2)$	Q<0.9	
1	А	338/342~(98%)	-0.13	7 (2%)	63	64	13, 20, 33, 67	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	1	GLN	4.5
1	А	338	ARG	4.1
1	А	145	ARG	2.5
1	А	278	TYR	2.3
1	А	141	GLU	2.1

#### 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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
6	EDO	А	408	4/4	0.64	0.17	40,44,45,53	0
5	ACT	А	407	4/4	0.72	0.28	26,29,31,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q<0.9
2	CL	А	401	1/1	0.84	0.07	48,48,48,48	0
4	CO2	А	406	3/3	0.84	0.12	43,43,47,52	0
6	EDO	А	409	4/4	0.88	0.23	34,38,38,42	0
6	EDO	А	411	4/4	0.89	0.33	34,34,34,35	0
7	PDO	А	412	5/5	0.89	0.20	28,29,34,34	0
8	PEG	А	414	7/7	0.89	0.13	30,33,37,37	0
2	CL	А	403	1/1	0.92	0.09	$47,\!47,\!47,\!47$	0
7	PDO	А	413	5/5	0.92	0.16	36, 38, 41, 41	0
2	CL	А	402	1/1	0.92	0.05	$51,\!51,\!51,\!51$	0
6	EDO	А	410	4/4	0.93	0.27	24,28,32,33	0
2	CL	А	404	1/1	0.95	0.13	49,49,49,49	0
3	SO2	А	405	3/3	0.96	0.05	34,34,48,51	0

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### 6.5 Other polymers (i)

There are no such residues in this entry.

