

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

#### Sep 24, 2023 – 09:28 AM EDT

PDB ID : 5UND

Title : Crystal Structure of CTCF(ZnF 4-10) With 28-mer DNA

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Deposited on : 2017-01-30

Resolution : 2.55 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 as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.35.1

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)

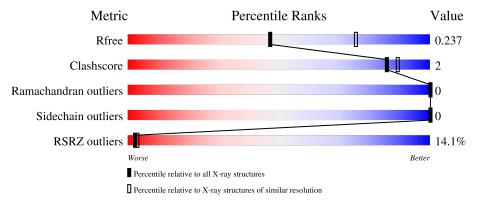
Validation Pipeline (wwPDB-VP) : 2.35.1

# 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 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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		
			8%		
1	A	202	81%	•	15%
			10%		
1	В	202	68%	28%	
			18%		
2	С	28	71%	21%	7%
			14%		
2	Е	28	82%	119	% 7%
			29%		
3	D	28	75%	18%	7%

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Mol	Chain	Length	Quality of chain		
			18%		
3	F	28	79%	11%	11%



# 2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 4723 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 Transcriptional repressor CTCF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171		C		0	S	0	0	0
			1338	829	259	232	18			
1	R	145	Total	$\mathbf{C}$	N	O	$\mathbf{S}$	0	9	0
1		140	1164	721	225	204	14		2	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	GLY	-	expression tag	UNP P49711
A	347	SER	-	expression tag	UNP P49711
В	346	GLY	-	expression tag	UNP P49711
В	347	SER	-	expression tag	UNP P49711

• Molecule 2 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	2 C	26	Total	С	N	О	Р	0	0	0
			521	248	91	157	25			
2	2 E	26	Total	С	N	О	Р	0	0	0
			521	248	91	157	25			U

• Molecule 3 is a DNA chain called DNA (26-MER).

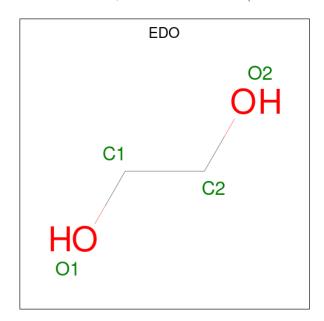
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	3 D	26	Total	С	N	О	Р	0	0	0
	20	541	252	111	152	26	U	0		
2	3 F	F 25	Total	С	N	О	Р	0	0	0
3			519	242	106	146	25	0	0	U

• Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	7	Total Zn 7 7	0	0
4	В	5	Total Zn 5 5	0	0

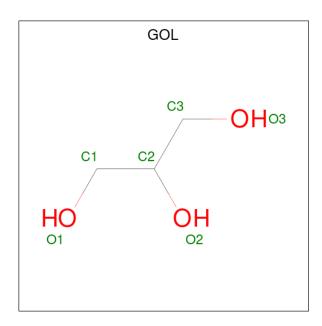
 $\bullet$  Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $\mathrm{C_2H_6O_2}).$ 



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	В	1	Total C O 4 2 2	0	0
5	Е	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0

 $\bullet$  Molecule 6 is GLYCEROL (three-letter code: GOL) (formula:  $\mathrm{C_3H_8O_3}).$ 





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

#### • Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	41	Total O 41 41	0	0
7	С	4	Total O 4 4	0	0
7	D	4	Total O 4 4	0	0
7	В	23	Total O 23 23	0	0
7	Е	1	Total O 1 1	0	0
7	F	4	Total O 4 4	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: Transcriptional repressor CTCF Chain A: 15% ASP MET HIS PHE LYS ARG TYR HIS • Molecule 1: Transcriptional repressor CTCF Chain B: 28% • Molecule 2: DNA (26-MER) Chain C: 71% 21% 7% • Molecule 2: DNA (26-MER) Chain E: 82% 11% • Molecule 3: DNA (26-MER) Chain D: 75% 18%



11%







# 4 Data and refinement statistics (i)

Property	Value	Source	
Space group	P 1 21 1	Depositor	
Cell constants	75.07Å 73.83Å 93.29Å	D	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $91.38^{\circ}$ $90.00^{\circ}$	Depositor	
Resolution (Å)	34.61 - 2.55	Depositor	
Resolution (A)	34.61 - 2.55	EDS	
% Data completeness	99.2 (34.61-2.55)	Depositor	
(in resolution range)	90.6 (34.61-2.55)	EDS	
$R_{merge}$	0.10	Depositor	
$R_{sum}$	(Not available)	Depositor	
$< I/\sigma(I) > 1$	1.20 (at 2.54Å)	Xtriage	
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor	
D D	0.211 , 0.237	Depositor	
$R, R_{free}$	0.212 , $0.237$	DCC	
$R_{free}$ test set	1669 reflections (5.02%)	wwPDB-VP	
Wilson B-factor (Å <sup>2</sup> )	54.7	Xtriage	
Anisotropy	0.400	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.27, 63.3	EDS	
L-test for twinning <sup>2</sup>	$< L > = 0.50, < L^2> = 0.33$	Xtriage	
	0.006 for k,h,-l		
Estimated twinning fraction	0.001  for -k,-h,-l	Xtriage	
	0.026  for h,-k,-l		
$F_o, F_c$ correlation	0.94	EDS	
Total number of atoms	4723	wwPDB-VP	
Average B, all atoms (Å <sup>2</sup> )	108.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.25% of the height of the origin peak. No significant pseudotranslation is detected.

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



<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: EDO, GOL, ZN

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		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.23	0/1372	0.38	0/1843	
1	В	0.23	0/1198	0.39	0/1609	
2	С	0.49	0/581	0.89	0/894	
2	Е	0.51	0/581	0.87	0/894	
3	D	0.48	0/609	0.74	0/939	
3	F	0.47	0/584	0.73	0/900	
All	All	0.38	0/4925	0.64	0/7079	

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	$\mathbf{H}(\mathbf{model})$	H(added)	Clashes	Symm-Clashes
1	A	1338	0	1210	5	0
1	В	1164	0	1066	5	0
2	С	521	0	290	3	0
2	Е	521	0	290	2	0
3	D	541	0	288	4	0
3	F	519	0	277	2	0
4	A	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	В	5	0	0	0	0
5	В	4	0	6	0	0
5	D	8	0	12	0	0
5	Е	4	0	6	0	0
5	F	8	0	12	0	0
6	В	6	0	8	0	0
7	A	41	0	0	0	0
7	В	23	0	0	0	0
7	С	4	0	0	0	0
7	D	4	0	0	0	0
7	Е	1	0	0	0	0
7	F	4	0	0	0	0
All	All	4723	0	3465	17	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 2.

All (17) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance}({ m \AA})$	$-$ overlap $(\AA)$
1:B:368:ARG:NH2	3:F:25:DG:N7	2.37	0.72
1:B:381:CYS:SG	1:B:397:HIS:CD2	2.93	0.61
2:E:1:DC:H2"	2:E:2:DG:C8	2.39	0.58
1:A:397:HIS:ND1	3:D:21:DA:OP1	2.36	0.50
1:A:356:CYS:SG	1:A:369:HIS:CE1	2.95	0.49
1:A:455:HIS:ND1	3:D:15:DG:OP1	2.43	0.49
1:A:371:ARG:HG3	1:A:376:GLU:HB3	1.96	0.48
3:D:21:DA:H2"	3:D:22:DG:H8	1.78	0.47
1:B:457:ARG:NH1	2:E:10:DC:OP2	2.48	0.47
1:B:437:PHE:HB3	1:B:452:LEU:HD22	1.97	0.46
2:C:18:DT:H2"	2:C:19:DG:C8	2.51	0.46
2:C:5:DC:H2'	2:C:6:DC:C6	2.51	0.45
1:B:362:GLU:HB2	1:B:365:LYS:HG2	2.00	0.44
3:F:27:DC:H2"	3:F:28:DG:C8	2.54	0.42
2:C:11:DT:H2"	2:C:12:DG:C8	2.54	0.41
3:D:27:DC:H2"	3:D:28:DG:C8	2.56	0.41
1:A:371:ARG:HD3	1:A:378:PRO:HD3	2.04	0.40

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	ntiles
1	A	169/202~(84%)	164 (97%)	5 (3%)	0	100	100
1	В	145/202~(72%)	142 (98%)	3 (2%)	0	100	100
All	All	314/404 (78%)	306 (98%)	8 (2%)	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 Outlier		Percentiles		
1	A	133/186 (72%)	133 (100%)	0	100	100	
1	В	120/186~(64%)	120 (100%)	0	100	100	
All	All	253/372~(68%)	253 (100%)	0	100	100	

There are no protein residues with a non-rotameric sidechain to report.

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 19 ligands modelled in this entry, 12 are monoatomic - leaving 7 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	Tuno	Chain	Res	Link	В	ond leng	gths	Bond angles		
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
5	EDO	D	102	-	3,3,3	0.46	0	2,2,2	0.33	0
5	EDO	Е	101	-	3,3,3	0.46	0	2,2,2	0.36	0
5	EDO	В	606	-	3,3,3	0.46	0	2,2,2	0.31	0
5	EDO	F	101	-	3,3,3	0.48	0	2,2,2	0.27	0
6	GOL	В	607	-	5,5,5	0.36	0	5,5,5	0.26	0
5	EDO	F	102	-	3,3,3	0.46	0	2,2,2	0.35	0
5	EDO	D	101	-	3,3,3	0.47	0	2,2,2	0.34	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	EDO	D	102	-	-	0/1/1/1	-
5	EDO	Е	101	-	-	0/1/1/1	-
5	EDO	В	606	-	-	0/1/1/1	-
5	EDO	F	101	-	-	0/1/1/1	-
6	GOL	В	607	-	-	2/4/4/4	-
5	EDO	F	102	-	-	0/1/1/1	-
5	EDO	D	101	-	-	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
6	В	607	GOL	O1-C1-C2-C3
6	В	607	GOL	O1-C1-C2-O2

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	<rsrz></rsrz>	# RSRZ > 2		2	$OWAB(\AA^2)$	Q < 0.9
1	A	171/202~(84%)	0.70	16 (9%)	8	10	37, 87, 157, 238	0
1	В	145/202 (71%)	0.83	21 (14%)	2	3	34, 75, 179, 241	0
2	С	$26/28 \; (92\%)$	0.44	5 (19%)	1	1	76, 98, 202, 260	0
2	E	$26/28 \; (92\%)$	1.00	4 (15%)	2	2	81, 111, 247, 316	0
3	D	26/28 (92%)	1.16	8 (30%)	0	0	54, 95, 207, 217	0
3	F	25/28 (89%)	1.14	5 (20%)	1	1	55, 112, 241, 261	0
All	All	419/516 (81%)	0.80	59 (14%)	2	3	34, 89, 202, 316	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	377	ARG	9.2
1	A	497	CYS	8.6
2	Е	24	DC	6.8
1	В	378	PRO	6.2
2	Е	25	DA	6.1
2	Е	26	DC	5.8
1	A	502	TYR	5.8
1	В	391	THR	5.8
1	В	392	TYR	5.2
3	F	7	DC	5.0
1	В	353	CYS	4.7
1	A	503	ALA	4.6
1	A	512	MET	4.6
1	В	375	GLY	4.5
1	В	389	ARG	4.3
1	В	358	TYR	4.2
3	F	6	DC	4.2
1	A	498	ASP	4.0
1	A	517	HIS	4.0

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Mol	Chain	$ hootnote{Res}$	Type	RSRZ
1	A	518	THR	4.0
1	A	504	CYS	4.0
1	В	359	ALA	3.9
3	D	3	DG	3.9
1	В	379	PHE	3.7
3	F	5	DG	3.7
1	В	376	GLU	3.6
1	A	499	GLN	3.6
2	С	26	DC	3.5
1	A	506	GLN	3.3
1	В	361	VAL	3.2
2	С	25	DA	3.2
1	В	354	SER	3.2
1	В	351	PHE	3.2
3	D	4	DT	3.0
3	D	9	DA	3.0
3	D	10	DC	3.0
3	F	8	DC	2.9
2	Е	23	DG	2.8
3	D	8	DC	2.6
1	В	355	MET	2.6
1	A	515	ARG	2.5
1	В	350	PRO	2.5
1	A	495	PHE	2.4
1	A	507	GLU	2.4
1	В	356	CYS	2.4
1	A	508	ARG	2.4
2	С	24	DC	2.4
1	A	383	LEU	2.4
3	D	7	DC	2.3
3	D	5	DG	2.3
1	В	445	VAL	2.2
1	В	374	THR	2.2
1	В	382	SER	2.1
2	С	23	DG	2.1
1	A	505	ARG	2.1
3	F	10	DC	2.1
2	С	22	DG	2.0
1	В	370	ILE	2.0
3	D	6	DC	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)

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	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
5	EDO	В	606	4/4	0.74	0.39	77,82,82,85	0
6	GOL	В	607	6/6	0.75	0.24	95,101,105,106	0
5	EDO	D	102	4/4	0.77	0.26	69,76,77,87	0
5	EDO	F	101	4/4	0.80	0.25	58,73,74,75	0
4	ZN	A	606	1/1	0.81	0.09	148,148,148,148	0
4	ZN	A	607	1/1	0.82	0.07	138,138,138,138	0
5	EDO	F	102	4/4	0.86	0.13	72,72,81,81	0
4	ZN	В	601	1/1	0.88	0.07	137,137,137,137	0
5	EDO	D	101	4/4	0.89	0.18	72,78,82,89	0
4	ZN	В	602	1/1	0.91	0.03	137,137,137,137	0
5	EDO	E	101	4/4	0.96	0.36	81,81,85,87	0
4	ZN	A	602	1/1	0.97	0.07	98,98,98,98	0
4	ZN	A	604	1/1	0.98	0.16	59,59,59,59	0
4	ZN	A	605	1/1	0.99	0.18	52,52,52,52	0
4	ZN	A	603	1/1	0.99	0.19	41,41,41,41	0
4	ZN	В	603	1/1	0.99	0.20	40,40,40,40	0
4	ZN	В	604	1/1	0.99	0.14	55,55,55,55	0
4	ZN	В	605	1/1	0.99	0.13	46,46,46,46	0
4	ZN	A	601	1/1	0.99	0.07	100,100,100,100	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

