



wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 09:12 am BST

PDB ID : 5ZMU
Title : Crystal structure of a cis-epoxysuccinate hydrolase producing D(-)-tartaric acids
Authors : Dong, S.; Liu, X.; Wang, X.; Feng, Y.
Deposited on : 2018-04-06
Resolution : 1.50 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
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)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

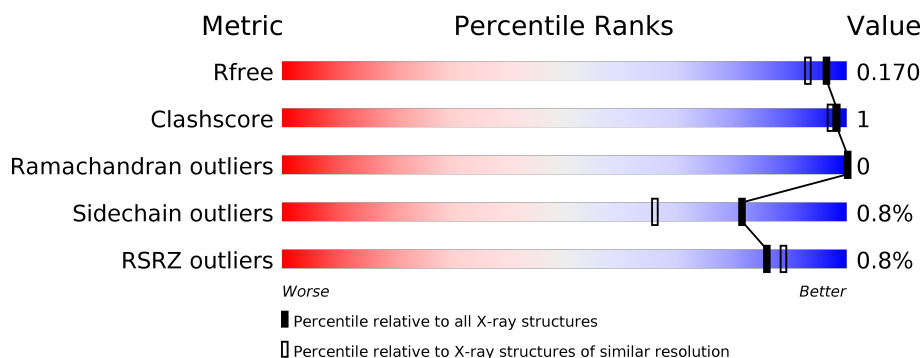
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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 $\leq 5\%$. 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	314	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 87%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 87% 11% </div> </div>
1	B	314	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 87%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 87% 11% </div> </div>
1	C	314	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 86%, yellow 1%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 86% 11% </div> </div>
1	D	314	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 89%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 89% 9% </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19369 atoms, of which 8854 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 Cis-epoxysuccinate hydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	281	Total	C	H	N	O	S	0	6	0
			4456	1411	2212	408	416	9			
1	B	281	Total	C	H	N	O	S	0	6	0
			4450	1410	2208	406	417	9			
1	C	280	Total	C	H	N	O	S	0	5	0
			4420	1403	2194	400	413	10			
1	D	287	Total	C	H	N	O	S	0	5	0
			4515	1429	2240	413	423	10			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP F1LJ99
A	2	GLY	-	expression tag	UNP F1LJ99
A	3	SER	-	expression tag	UNP F1LJ99
A	4	SER	-	expression tag	UNP F1LJ99
A	5	HIS	-	expression tag	UNP F1LJ99
A	6	HIS	-	expression tag	UNP F1LJ99
A	7	HIS	-	expression tag	UNP F1LJ99
A	8	HIS	-	expression tag	UNP F1LJ99
A	9	HIS	-	expression tag	UNP F1LJ99
A	10	HIS	-	expression tag	UNP F1LJ99
A	11	SER	-	expression tag	UNP F1LJ99
A	12	SER	-	expression tag	UNP F1LJ99
A	13	GLY	-	expression tag	UNP F1LJ99
A	14	LEU	-	expression tag	UNP F1LJ99
A	15	VAL	-	expression tag	UNP F1LJ99
A	16	PRO	-	expression tag	UNP F1LJ99
A	17	ARG	-	expression tag	UNP F1LJ99
A	18	GLY	-	expression tag	UNP F1LJ99
A	19	SER	-	expression tag	UNP F1LJ99
A	20	HIS	-	expression tag	UNP F1LJ99
B	1	MET	-	expression tag	UNP F1LJ99

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Chain	Residue	Modelled	Actual	Comment	Reference
B	2	GLY	-	expression tag	UNP F1LJ99
B	3	SER	-	expression tag	UNP F1LJ99
B	4	SER	-	expression tag	UNP F1LJ99
B	5	HIS	-	expression tag	UNP F1LJ99
B	6	HIS	-	expression tag	UNP F1LJ99
B	7	HIS	-	expression tag	UNP F1LJ99
B	8	HIS	-	expression tag	UNP F1LJ99
B	9	HIS	-	expression tag	UNP F1LJ99
B	10	HIS	-	expression tag	UNP F1LJ99
B	11	SER	-	expression tag	UNP F1LJ99
B	12	SER	-	expression tag	UNP F1LJ99
B	13	GLY	-	expression tag	UNP F1LJ99
B	14	LEU	-	expression tag	UNP F1LJ99
B	15	VAL	-	expression tag	UNP F1LJ99
B	16	PRO	-	expression tag	UNP F1LJ99
B	17	ARG	-	expression tag	UNP F1LJ99
B	18	GLY	-	expression tag	UNP F1LJ99
B	19	SER	-	expression tag	UNP F1LJ99
B	20	HIS	-	expression tag	UNP F1LJ99
C	1	MET	-	expression tag	UNP F1LJ99
C	2	GLY	-	expression tag	UNP F1LJ99
C	3	SER	-	expression tag	UNP F1LJ99
C	4	SER	-	expression tag	UNP F1LJ99
C	5	HIS	-	expression tag	UNP F1LJ99
C	6	HIS	-	expression tag	UNP F1LJ99
C	7	HIS	-	expression tag	UNP F1LJ99
C	8	HIS	-	expression tag	UNP F1LJ99
C	9	HIS	-	expression tag	UNP F1LJ99
C	10	HIS	-	expression tag	UNP F1LJ99
C	11	SER	-	expression tag	UNP F1LJ99
C	12	SER	-	expression tag	UNP F1LJ99
C	13	GLY	-	expression tag	UNP F1LJ99
C	14	LEU	-	expression tag	UNP F1LJ99
C	15	VAL	-	expression tag	UNP F1LJ99
C	16	PRO	-	expression tag	UNP F1LJ99
C	17	ARG	-	expression tag	UNP F1LJ99
C	18	GLY	-	expression tag	UNP F1LJ99
C	19	SER	-	expression tag	UNP F1LJ99
C	20	HIS	-	expression tag	UNP F1LJ99
D	1	MET	-	expression tag	UNP F1LJ99
D	2	GLY	-	expression tag	UNP F1LJ99
D	3	SER	-	expression tag	UNP F1LJ99

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Chain	Residue	Modelled	Actual	Comment	Reference
D	4	SER	-	expression tag	UNP F1LJ99
D	5	HIS	-	expression tag	UNP F1LJ99
D	6	HIS	-	expression tag	UNP F1LJ99
D	7	HIS	-	expression tag	UNP F1LJ99
D	8	HIS	-	expression tag	UNP F1LJ99
D	9	HIS	-	expression tag	UNP F1LJ99
D	10	HIS	-	expression tag	UNP F1LJ99
D	11	SER	-	expression tag	UNP F1LJ99
D	12	SER	-	expression tag	UNP F1LJ99
D	13	GLY	-	expression tag	UNP F1LJ99
D	14	LEU	-	expression tag	UNP F1LJ99
D	15	VAL	-	expression tag	UNP F1LJ99
D	16	PRO	-	expression tag	UNP F1LJ99
D	17	ARG	-	expression tag	UNP F1LJ99
D	18	GLY	-	expression tag	UNP F1LJ99
D	19	SER	-	expression tag	UNP F1LJ99
D	20	HIS	-	expression tag	UNP F1LJ99

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0

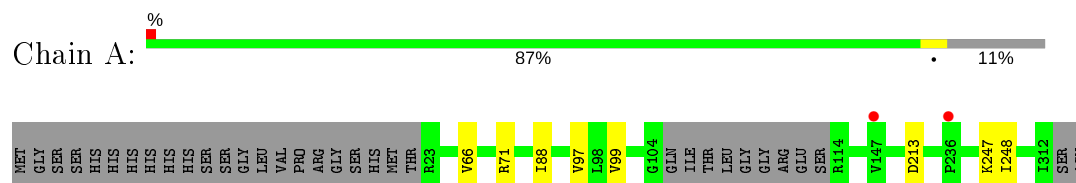
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	392	Total O 392 392	0	0
3	B	357	Total O 357 357	0	0
3	C	415	Total O 415 415	0	0
3	D	360	Total O 360 360	0	0

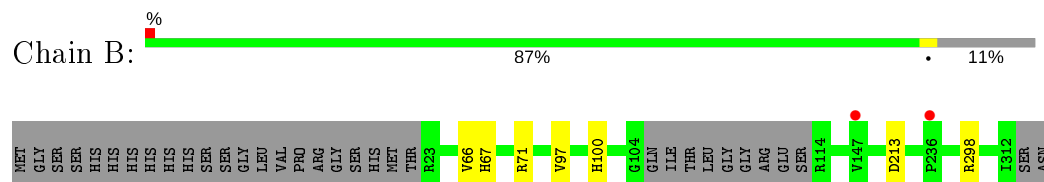
3 Residue-property plots [i](#)

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.

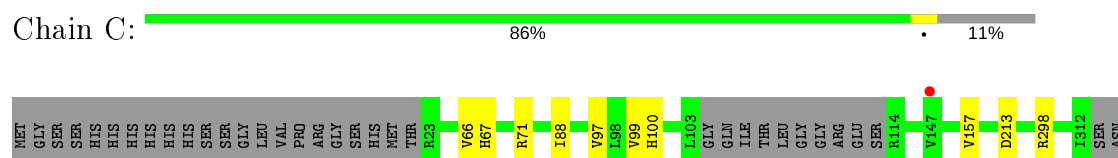
- Molecule 1: Cis-epoxysuccinate hydrolase



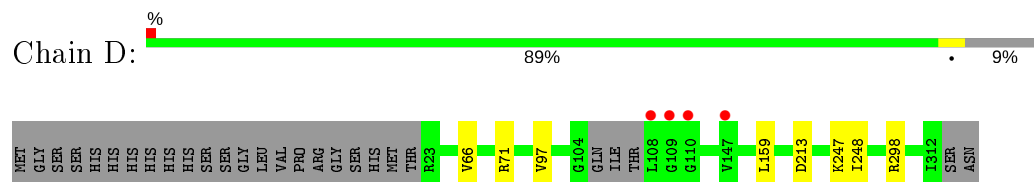
- Molecule 1: Cis-epoxysuccinate hydrolase



- Molecule 1: Cis-epoxysuccinate hydrolase



- Molecule 1: Cis-epoxysuccinate hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	78.49Å 94.25Å 94.36Å 90.00° 95.90° 90.00°	Depositor
Resolution (Å)	40.35 – 1.50 40.35 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.35-1.50) 93.9 (40.35-1.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 1.50Å)	Xtriage
Refinement program	phenix.refine 1.13_2998, PHENIX 1.13_2998	Depositor
R, R_{free}	0.149 , 0.170 0.149 , 0.170	Depositor DCC
R_{free} test set	2000 reflections (0.92%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19369	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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: 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	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.53	0/2295	0.67	1/3117 (0.0%)
1	B	0.55	0/2293	0.69	1/3115 (0.0%)
1	C	0.54	0/2277	0.70	1/3094 (0.0%)
1	D	0.55	0/2326	0.69	1/3157 (0.0%)
All	All	0.54	0/9191	0.69	4/12483 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	71	ARG	NE-CZ-NH1	8.08	124.34	120.30
1	D	71	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	B	71	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	A	71	ARG	NE-CZ-NH1	5.83	123.21	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	A	2244	2212	2207	3	0
1	B	2242	2208	2202	2	0
1	C	2226	2194	2189	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2275	2240	2235	3	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	392	0	0	0	1
3	B	357	0	0	0	1
3	C	415	0	0	0	0
3	D	360	0	0	0	0
All	All	10515	8854	8833	11	1

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 1.

The worst 5 of 11 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:C:66[B]:VAL:HG23	1:C:97:VAL:HG21	1.72	0.70
1:A:66[B]:VAL:HG23	1:A:97:VAL:HG21	1.74	0.68
1:B:66[B]:VAL:HG23	1:B:97:VAL:HG21	1.73	0.68
1:D:66[B]:VAL:HG23	1:D:97:VAL:HG21	1.87	0.56
1:A:88:ILE:HG12	1:A:99:VAL:HG11	1.97	0.46

All (1) 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)
3:A:870:HOH:O	3:B:843:HOH:O[2_655]	2.10	0.10

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	A	283/314 (90%)	279 (99%)	4 (1%)	0	100	100
1	B	283/314 (90%)	280 (99%)	3 (1%)	0	100	100
1	C	281/314 (90%)	277 (99%)	4 (1%)	0	100	100
1	D	288/314 (92%)	284 (99%)	4 (1%)	0	100	100
All	All	1135/1256 (90%)	1120 (99%)	15 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	A	233/255 (91%)	232 (100%)	1 (0%)	91	82
1	B	233/255 (91%)	231 (99%)	2 (1%)	78	61
1	C	232/255 (91%)	230 (99%)	2 (1%)	78	61
1	D	236/255 (92%)	234 (99%)	2 (1%)	81	66
All	All	934/1020 (92%)	927 (99%)	7 (1%)	81	69

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

Mol	Chain	Res	Type
1	C	213	ASP
1	D	298	ARG
1	C	298	ARG
1	B	213	ASP
1	D	213	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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	A	281/314 (89%)	-0.50	2 (0%) 87 90	11, 17, 34, 47	0
1	B	281/314 (89%)	-0.49	2 (0%) 87 90	11, 17, 33, 46	0
1	C	280/314 (89%)	-0.52	1 (0%) 92 94	11, 16, 31, 46	0
1	D	287/314 (91%)	-0.53	4 (1%) 75 79	11, 16, 34, 61	0
All	All	1129/1256 (89%)	-0.51	9 (0%) 86 89	11, 16, 33, 61	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	147	VAL	3.8
1	D	110	GLY	3.5
1	A	147	VAL	2.8
1	B	236	PRO	2.6
1	D	109	GLY	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](#)

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, 95th 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(\AA^2)	Q<0.9
2	ZN	C	400	1/1	0.99	0.04	16,16,16,16	1
2	ZN	B	400	1/1	0.99	0.04	15,15,15,15	1
2	ZN	A	400	1/1	1.00	0.04	14,14,14,14	1
2	ZN	D	400	1/1	1.00	0.04	15,15,15,15	1

6.5 Other polymers [i](#)

There are no such residues in this entry.