



# wwPDB X-ray Structure Validation Summary Report i

Feb 27, 2014 – 11:12 PM GMT

PDB ID : 1J5S  
Title : Crystal structure of uronate isomerase (TM0064) from *Thermotoga maritima*  
at 2.85 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : 2002-07-02  
Resolution : 2.85 Å (reported)

This is a wwPDB validation summary report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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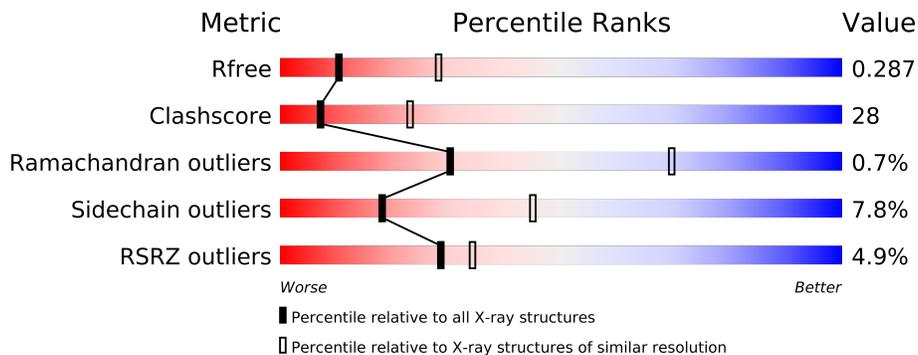
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.85 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))
$R_{free}$	66092	1524 (2.90-2.82)
Clashscore	79885	1879 (2.90-2.82)
Ramachandran outliers	78287	1824 (2.90-2.82)
Sidechain outliers	78261	1827 (2.90-2.82)
RSRZ outliers	66119	1526 (2.90-2.82)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	463	
1	B	463	
1	C	463	

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 11175 atoms, of which 0 are hydrogen and 0 are deuterium.

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 URONATE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	451	Total 3696	C 2367	N 632	O 680	S 17	81	0	0
1	B	450	Total 3686	C 2361	N 629	O 679	S 17	86	0	0
1	C	450	Total 3686	C 2361	N 629	O 679	S 17	83	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	LEADER SEQUENCE	UNP Q9WXR9
A	-10	GLY	-	LEADER SEQUENCE	UNP Q9WXR9
A	-9	SER	-	LEADER SEQUENCE	UNP Q9WXR9
A	-8	ASP	-	LEADER SEQUENCE	UNP Q9WXR9
A	-7	LYS	-	LEADER SEQUENCE	UNP Q9WXR9
A	-6	ILE	-	LEADER SEQUENCE	UNP Q9WXR9
A	-5	HIS	-	EXPRESSION TAG	UNP Q9WXR9
A	-4	HIS	-	EXPRESSION TAG	UNP Q9WXR9
A	-3	HIS	-	EXPRESSION TAG	UNP Q9WXR9
A	-2	HIS	-	EXPRESSION TAG	UNP Q9WXR9
A	-1	HIS	-	EXPRESSION TAG	UNP Q9WXR9
A	0	HIS	-	EXPRESSION TAG	UNP Q9WXR9
B	-11	MET	-	LEADER SEQUENCE	UNP Q9WXR9
B	-10	GLY	-	LEADER SEQUENCE	UNP Q9WXR9
B	-9	SER	-	LEADER SEQUENCE	UNP Q9WXR9
B	-8	ASP	-	LEADER SEQUENCE	UNP Q9WXR9
B	-7	LYS	-	LEADER SEQUENCE	UNP Q9WXR9
B	-6	ILE	-	LEADER SEQUENCE	UNP Q9WXR9
B	-5	HIS	-	EXPRESSION TAG	UNP Q9WXR9
B	-4	HIS	-	EXPRESSION TAG	UNP Q9WXR9
B	-3	HIS	-	EXPRESSION TAG	UNP Q9WXR9
B	-2	HIS	-	EXPRESSION TAG	UNP Q9WXR9
B	-1	HIS	-	EXPRESSION TAG	UNP Q9WXR9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	-11	MET	-	LEADER SEQUENCE	UNP Q9WXR9
C	-10	GLY	-	LEADER SEQUENCE	UNP Q9WXR9
C	-9	SER	-	LEADER SEQUENCE	UNP Q9WXR9
C	-8	ASP	-	LEADER SEQUENCE	UNP Q9WXR9
C	-7	LYS	-	LEADER SEQUENCE	UNP Q9WXR9
C	-6	ILE	-	LEADER SEQUENCE	UNP Q9WXR9
C	-5	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	-4	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	-3	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	-2	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	-1	HIS	-	EXPRESSION TAG	UNP Q9WXR9
C	0	HIS	-	EXPRESSION TAG	UNP Q9WXR9

- Molecule 2 is water.

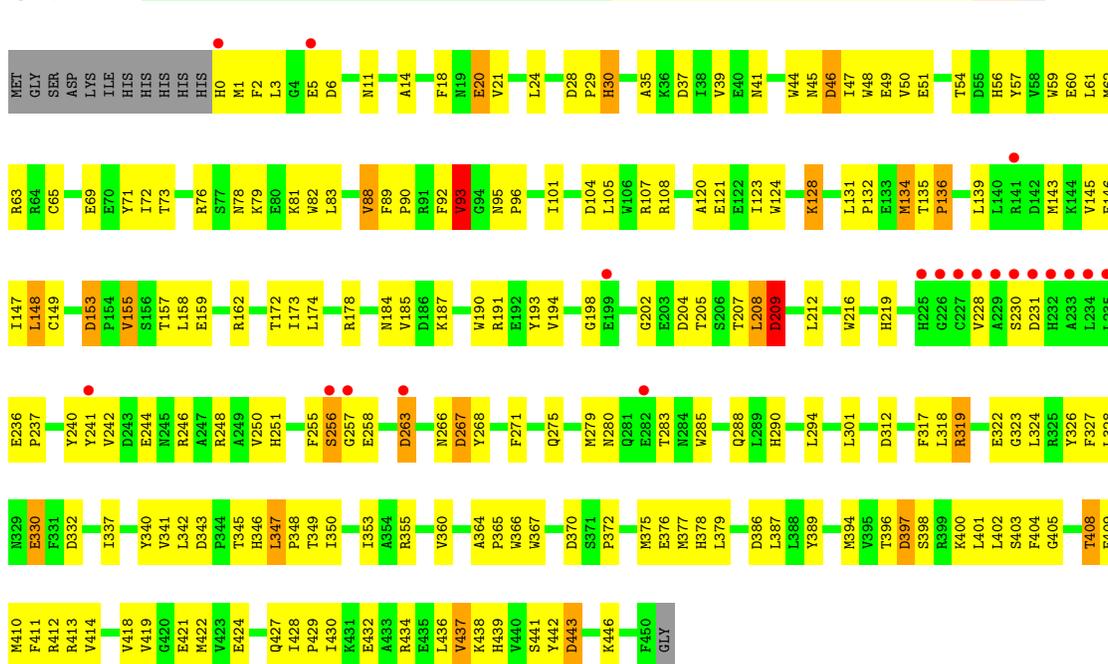
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	40	Total O 40 40	0	0
2	B	33	Total O 33 33	0	0
2	C	34	Total O 34 34	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

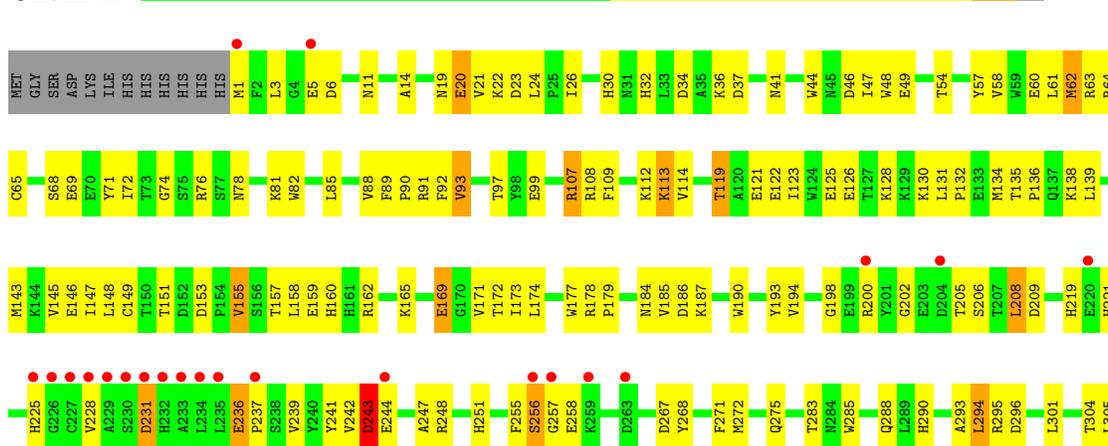
- Molecule 1: URONATE ISOMERASE

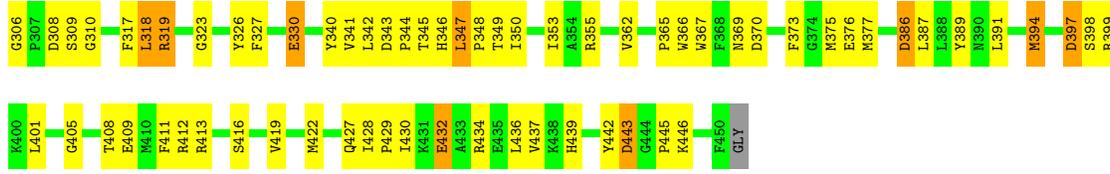
Chain A:



- Molecule 1: URONATE ISOMERASE

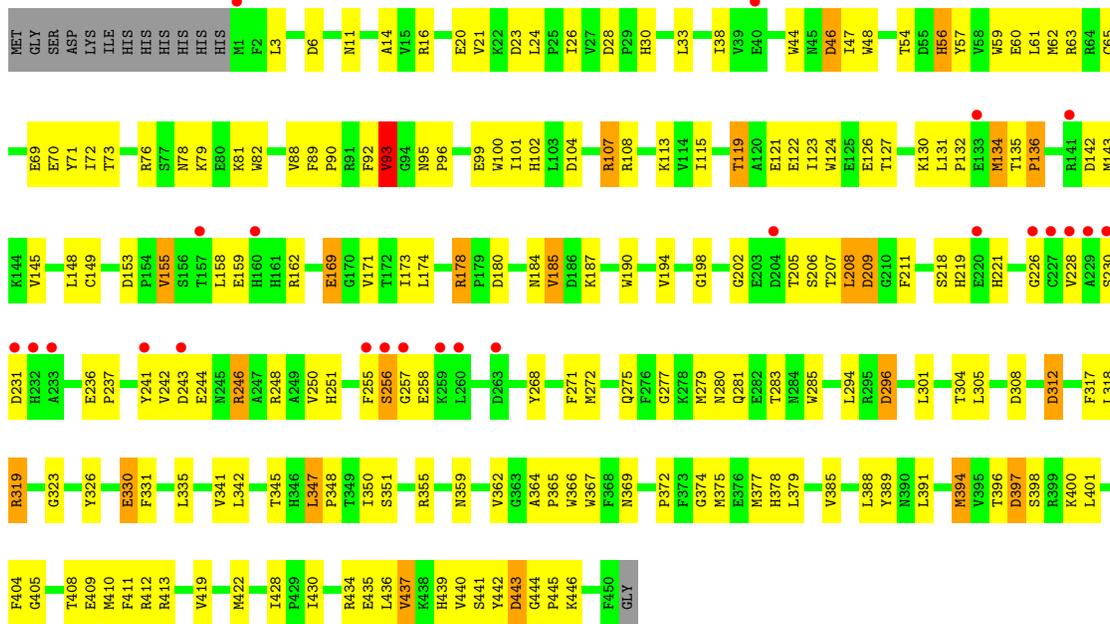
Chain B:





• Molecule 1: URONATE ISOMERASE

Chain C:



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.41Å 79.96Å 89.43Å 115.73° 97.57° 110.44°	Depositor
Resolution (Å)	24.99 – 2.85 24.99 – 2.75	Depositor EDS
% Data completeness (in resolution range)	96.6 (24.99-2.85) 96.8 (24.99-2.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.76Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.234 , 0.274 0.250 , 0.287	Depositor DCC
$R_{free}$ test set	2153 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.1	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 38.6	EDS
Estimated twinning fraction	0.021 for k,h,-h-k-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 84993 reflections	Xtriage
$F_o, F_c$ correlation	0.85	EDS
Total number of atoms	11175	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality (i)

### 5.1 Standard geometry (i)

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	0.52	0/3793	0.88	12/5136 (0.2%)
1	B	0.54	0/3782	0.90	12/5121 (0.2%)
1	C	0.54	0/3782	0.89	12/5121 (0.2%)
All	All	0.54	0/11357	0.89	36/15378 (0.2%)

There are no bond length outliers.

The worst 5 of 36 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	243	ASP	N-CA-CB	-9.25	93.94	110.60
1	B	243	ASP	CB-CG-OD2	8.96	126.37	118.30
1	A	104	ASP	CB-CG-OD2	7.03	124.62	118.30
1	A	132	PRO	CA-N-CD	-6.98	101.72	111.50
1	B	23	ASP	CB-CG-OD2	6.65	124.29	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3696	0	3621	217	7
1	B	3686	0	3614	202	11
1	C	3686	0	3614	223	4

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	40	0	0	3	0
2	B	33	0	0	4	0
2	C	34	0	0	5	0
All	All	11175	0	10849	603	11

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 28.

The worst 5 of 603 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:194:VAL:HG12	1:A:205:THR:CG2	1.53	1.34
1:A:251:HIS:ND1	2:A:454:HOH:O	1.65	1.21
1:B:62:MET:CE	1:B:72:ILE:HG12	1.73	1.19
1:A:194:VAL:CG1	1:A:205:THR:CG2	2.22	1.17
1:C:400:LYS:HD3	2:C:458:HOH:O	1.40	1.17

The worst 5 of 11 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	Distance(Å)	Clash(Å)
1:A:204:ASP:OD2	1:B:112:LYS:NZ[1_655]	0.84	1.36
1:A:128:LYS:NZ	1:B:169:GLU:OE2[1_665]	1.20	1.00
1:B:128:LYS:NZ	1:C:169:GLU:CD[1_556]	1.49	0.71
1:B:128:LYS:NZ	1:C:169:GLU:OE2[1_556]	1.51	0.69
1:A:255:PHE:O	1:B:5:GLU:OE2[1_655]	1.58	0.62

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	449/463 (97%)	425 (95%)	20 (4%)	4 (1%)	25 62
1	B	448/463 (97%)	420 (94%)	25 (6%)	3 (1%)	30 70
1	C	448/463 (97%)	422 (94%)	23 (5%)	3 (1%)	30 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1345/1389 (97%)	1267 (94%)	68 (5%)	10 (1%)	30 70

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	257	GLY
1	B	257	GLY
1	C	257	GLY
1	B	169	GLU
1	B	243	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	A	399/409 (98%)	367 (92%)	32 (8%)	17 44
1	B	398/409 (97%)	367 (92%)	31 (8%)	18 45
1	C	398/409 (97%)	368 (92%)	30 (8%)	19 48
All	All	1195/1227 (97%)	1102 (92%)	93 (8%)	18 45

5 of 93 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	155	VAL
1	B	330	GLU
1	C	330	GLU
1	B	171	VAL
1	B	256	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	288	GLN
1	C	56	HIS
1	C	251	HIS

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Mol	Chain	Res	Type
1	B	221	HIS
1	C	288	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

### 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

### 6.1 Protein, DNA and RNA chains

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	451/463 (97%)	-0.13	20 (4%) 33 39	8, 28, 84, 122	18 (3%)
1	B	450/463 (97%)	-0.10	22 (4%) 28 34	4, 27, 83, 122	19 (4%)
1	C	450/463 (97%)	-0.11	24 (5%) 25 31	8, 28, 83, 122	19 (4%)
All	All	1351/1389 (97%)	-0.11	66 (4%) 28 34	4, 28, 84, 122	56 (4%)

The worst 5 of 66 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	ALA	5.4
1	C	229	ALA	5.3
1	A	232	HIS	5.2
1	B	229	ALA	5.0
1	A	227	CYS	5.0

### 6.2 Non-standard residues in protein, DNA, RNA chains

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

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.