



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:26 am GMT

PDB ID : 1V5X  
Title : Crystal structure of Phosphoribosyl anthranilate isomerase from *Thermus Thermophilus*  
Authors : Taka, J.; Kunishima, N.; Yutani, K.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2003-11-26  
Resolution : 2.00 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

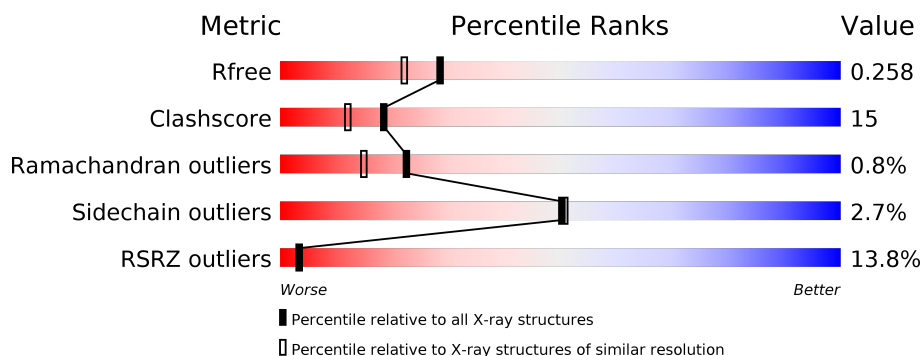
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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. 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	203	<div> <div>26%</div> <div>61%</div> <div>34%</div> <div>..</div> </div>
1	B	203	<div> <div>%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3357 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 Phosphoribosylanthranilate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	4	0	0
			1528	981	276	268	3			
1	B	200	Total	C	N	O	S	0	0	0
			1528	981	276	268	3			

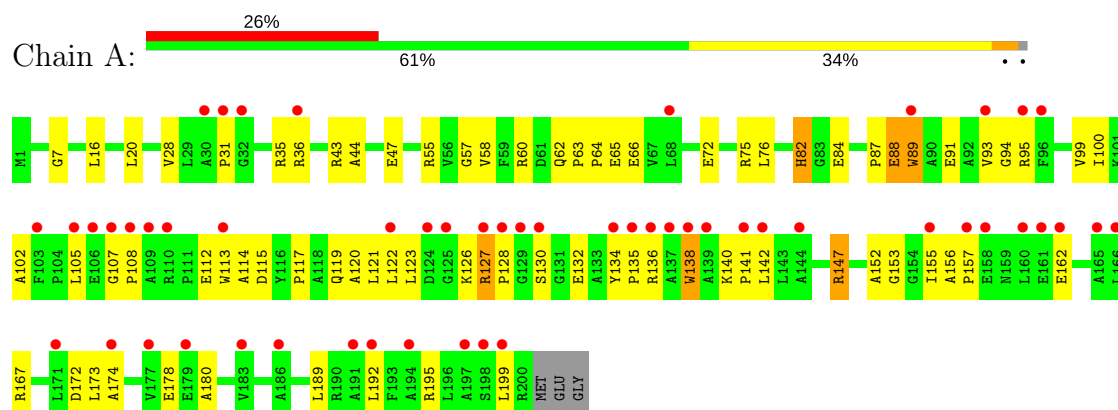
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	104	Total	O	0	0
			104	104		
2	B	197	Total	O	0	0
			197	197		

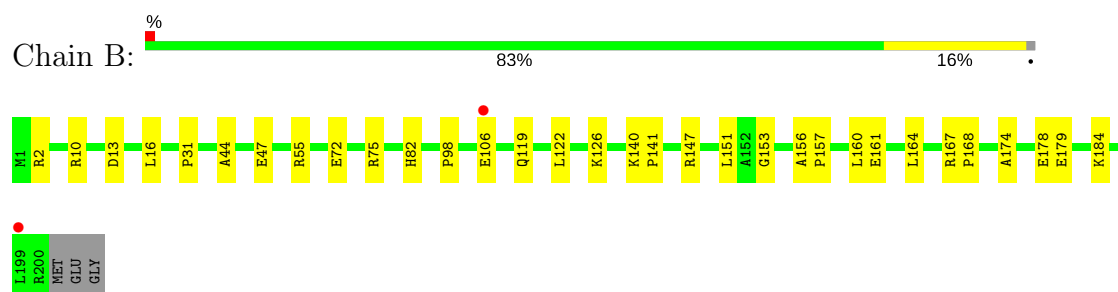
### 3 Residue-property plots

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



#### • Molecule 1: Phosphoribosylanthranilate isomerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	144.21 Å   144.21 Å   144.21 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	35.00 – 2.00 34.97 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.00-2.00) 99.8 (34.97-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.87 (at 2.00 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.262   ,   0.296 0.264   ,   0.258	Depositor DCC
$R_{free}$ test set	1761 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 61.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3357	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<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.

## 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	0.43	1/1563 (0.1%)	0.64	0/2123
1	B	0.37	0/1563	0.62	0/2123
All	All	0.40	1/3126 (0.0%)	0.63	0/4246

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	TRP	NE1-CE2	8.73	1.48	1.37

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	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1528	0	1572	72	0
1	B	1528	0	1572	20	0
2	A	104	0	0	5	0
2	B	197	0	0	1	0
All	All	3357	0	3144	92	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 15.

All (92) 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:147:ARG:HB3	1:A:147:ARG:HH11	1.37	0.87
1:A:136:ARG:HG3	1:A:162:GLU:HB3	1.67	0.77
1:B:174:ALA:O	1:B:178:GLU:HG3	1.86	0.75
1:A:87:PRO:HG2	1:A:88:GLU:OE1	1.86	0.74
1:A:60:ARG:HH22	1:A:128:PRO:HG3	1.51	0.74
1:B:10:ARG:HD3	1:B:13:ASP:OD2	1.87	0.74
1:A:174:ALA:O	1:A:178:GLU:HG3	1.88	0.73
1:A:152:ALA:HA	1:A:155:ILE:CD1	2.23	0.69
1:A:157:PRO:HB3	1:A:192:LEU:HA	1.74	0.67
1:A:112:GLU:O	1:A:115:ASP:HB2	1.93	0.67
1:A:127:ARG:HB2	1:A:130:SER:OG	1.97	0.64
1:A:134:TYR:HB2	1:A:135:PRO:HD2	1.78	0.64
1:B:98:PRO:HA	1:B:119:GLN:NE2	2.12	0.64
1:A:152:ALA:HA	1:A:155:ILE:HD12	1.80	0.64
1:A:152:ALA:CA	1:A:155:ILE:HD11	2.28	0.63
1:A:105:LEU:HD11	1:A:123:LEU:HB3	1.79	0.63
1:A:126:LYS:HB3	1:A:127:ARG:HD3	1.81	0.62
1:A:167:ARG:HH11	1:A:167:ARG:HB3	1.63	0.62
1:B:178:GLU:HG2	1:B:184:LYS:HD3	1.80	0.62
1:A:31:PRO:HA	1:A:36:ARG:HH11	1.65	0.62
1:A:119:GLN:HG3	2:A:243:HOH:O	2.00	0.61
1:A:147:ARG:NH1	1:A:147:ARG:HB3	2.15	0.60
1:A:195:ARG:O	1:A:199:LEU:HG	2.02	0.60
1:B:55:ARG:HH11	1:B:55:ARG:CG	2.16	0.58
1:A:122:LEU:HD21	1:A:152:ALA:CB	2.33	0.58
1:B:31:PRO:HD2	2:B:323:HOH:O	2.04	0.58
1:A:112:GLU:HB2	2:A:291:HOH:O	2.05	0.56
1:A:100:ILE:HG12	1:A:120:ALA:HB3	1.88	0.56
1:A:64:PRO:HD3	1:A:89:TRP:CD2	2.41	0.55
1:A:127:ARG:HH21	1:A:132:GLU:HG2	1.70	0.55
1:A:55:ARG:HH11	1:A:55:ARG:HG3	1.72	0.55
1:A:127:ARG:HH11	1:A:127:ARG:HG2	1.72	0.55
1:A:31:PRO:HA	1:A:36:ARG:NH1	2.20	0.55
1:A:44:ALA:HA	1:A:47:GLU:OE2	2.07	0.54
1:A:113:TRP:C	1:A:115:ASP:H	2.10	0.54
1:A:173:LEU:HD22	1:A:192:LEU:HD22	1.89	0.54
1:A:82:HIS:HA	1:A:102:ALA:HB3	1.89	0.53
1:A:123:LEU:HD21	1:A:142:LEU:CD2	2.38	0.53
1:A:63:PRO:HB2	1:A:65:GLU:HG2	1.90	0.53
1:A:156:ALA:HB1	1:A:157:PRO:HD2	1.91	0.52
1:A:114:ALA:HB1	1:A:147:ARG:HG3	1.92	0.52
1:A:91:GLU:OE2	1:A:117:PRO:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ARG:HB3	1:A:167:ARG:NH1	2.25	0.52
1:A:123:LEU:HD21	1:A:142:LEU:HD23	1.92	0.52
1:A:152:ALA:CA	1:A:155:ILE:CD1	2.88	0.52
1:A:178:GLU:HB3	1:A:180:ALA:O	2.10	0.51
1:A:173:LEU:HD22	1:A:192:LEU:CD2	2.41	0.51
1:A:108:PRO:HD2	2:A:290:HOH:O	2.11	0.51
1:A:127:ARG:HH21	1:A:132:GLU:CG	2.24	0.50
1:B:106:GLU:OE1	1:B:126:LYS:HE2	2.12	0.50
1:A:63:PRO:O	1:A:66:GLU:HB3	2.11	0.50
1:A:91:GLU:HB3	1:A:95:ARG:NH2	2.27	0.50
1:A:113:TRP:C	1:A:115:ASP:N	2.65	0.50
1:A:122:LEU:HD21	1:A:152:ALA:HB2	1.93	0.50
1:B:167:ARG:HH11	1:B:167:ARG:HG3	1.77	0.50
1:A:114:ALA:HA	1:A:121:LEU:HD11	1.94	0.49
1:A:84:GLU:OE1	1:A:84:GLU:HA	2.12	0.49
1:A:140:LYS:HA	2:A:301:HOH:O	2.13	0.48
1:B:72:GLU:O	1:B:75:ARG:NH1	2.47	0.48
1:A:43:ARG:NH1	2:A:231:HOH:O	2.45	0.47
1:A:122:LEU:HD21	1:A:152:ALA:HB3	1.96	0.47
1:B:160:LEU:O	1:B:164:LEU:HB2	2.15	0.46
1:A:72:GLU:O	1:A:75:ARG:NH1	2.47	0.46
1:A:127:ARG:N	1:A:127:ARG:HD3	2.31	0.46
1:A:93:VAL:HG23	1:A:99:VAL:HG21	1.97	0.45
1:A:7:GLY:HA3	1:A:35:ARG:HD3	1.98	0.45
1:A:28:VAL:HA	1:A:58:VAL:HB	1.98	0.44
1:A:36:ARG:HD2	1:A:36:ARG:HH11	1.61	0.44
1:A:155:ILE:HD12	1:A:172:ASP:O	2.18	0.44
1:A:93:VAL:HG23	1:A:94:GLY:N	2.33	0.44
1:B:55:ARG:NH1	1:B:55:ARG:CG	2.77	0.43
1:A:107:GLY:HA2	1:A:138:TRP:CZ2	2.53	0.43
1:B:44:ALA:HA	1:B:47:GLU:OE2	2.18	0.43
1:A:93:VAL:C	1:A:95:ARG:H	2.22	0.43
1:A:127:ARG:HG2	1:A:127:ARG:NH1	2.34	0.43
1:A:57:GLY:CA	1:A:76:LEU:HD13	2.48	0.42
1:B:161:GLU:CD	1:B:161:GLU:H	2.22	0.42
1:A:20:LEU:CD1	1:A:189:LEU:HB3	2.49	0.42
1:B:122:LEU:C	1:B:122:LEU:HD23	2.41	0.41
1:A:136:ARG:CG	1:A:162:GLU:HB3	2.43	0.41
1:B:2:ARG:HG3	1:B:2:ARG:NH1	2.36	0.41
1:A:140:LYS:N	1:A:141:PRO:HD2	2.36	0.41
1:A:63:PRO:HA	1:A:64:PRO:HD3	1.96	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLN:HB3	1:A:66:GLU:HG2	2.03	0.41
1:A:107:GLY:HA2	1:A:138:TRP:CE2	2.56	0.40
1:B:55:ARG:NH1	1:B:55:ARG:HG3	2.35	0.40
1:B:140:LYS:N	1:B:141:PRO:HD2	2.36	0.40
1:B:147:ARG:HD2	1:B:147:ARG:HH11	1.63	0.40
1:A:63:PRO:HG2	1:A:66:GLU:HB2	2.03	0.40
1:A:88:GLU:H	1:A:88:GLU:CD	2.25	0.40
1:B:151:LEU:HD12	1:B:168:PRO:HG3	2.02	0.40
1:B:156:ALA:HB1	1:B:157:PRO:CD	2.52	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	Percentiles	
1	A	198/203 (98%)	170 (86%)	26 (13%)	2 (1%)	18	10
1	B	198/203 (98%)	192 (97%)	5 (2%)	1 (0%)	32	26
All	All	396/406 (98%)	362 (91%)	31 (8%)	3 (1%)	22	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	TRP
1	A	153	GLY
1	B	153	GLY

### 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	148/150 (99%)	143 (97%)	5 (3%)	42	40
1	B	148/150 (99%)	145 (98%)	3 (2%)	60	64
All	All	296/300 (99%)	288 (97%)	8 (3%)	50	51

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	82	HIS
1	A	88	GLU
1	A	127	ARG
1	A	147	ARG
1	B	16	LEU
1	B	82	HIS
1	B	179	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	200/203 (98%)	1.46	53 (26%) 1 1	19, 56, 85, 89	29 (14%)
1	B	200/203 (98%)	0.12	2 (1%) 82 82	17, 30, 50, 65	4 (2%)
All	All	400/406 (98%)	0.79	55 (13%) 3 3	17, 40, 81, 89	33 (8%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	127	ARG	6.2
1	A	165	ALA	5.9
1	A	105	LEU	5.6
1	A	137	ALA	5.6
1	A	113	TRP	5.3
1	A	198	SER	5.0
1	A	138	TRP	4.6
1	A	89	TRP	4.5
1	A	128	PRO	4.5
1	A	139	ALA	4.3
1	A	129	GLY	4.1
1	A	160	LEU	4.1
1	A	135	PRO	3.7
1	A	134	TYR	3.7
1	A	197	ALA	3.6
1	A	141	PRO	3.5
1	A	109	ALA	3.5
1	A	108	PRO	3.3
1	A	125	GLY	3.3
1	A	124	ASP	3.2
1	A	103	PHE	3.2
1	A	110	ARG	3.2
1	A	191	ALA	3.2
1	A	177	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	199	LEU	3.1
1	A	162	GLU	3.1
1	A	106	GLU	3.1
1	A	142	LEU	3.0
1	A	158	GLU	3.0
1	A	130	SER	2.8
1	A	136	ARG	2.8
1	A	30	ALA	2.7
1	A	174	ALA	2.7
1	A	183	VAL	2.7
1	A	161	GLU	2.7
1	A	179	GLU	2.6
1	A	192	LEU	2.6
1	A	122	LEU	2.6
1	A	171	LEU	2.6
1	A	157	PRO	2.6
1	A	194	ALA	2.5
1	A	95	ARG	2.5
1	A	166	LEU	2.5
1	B	199	LEU	2.4
1	A	93	VAL	2.4
1	A	68	LEU	2.3
1	A	32	GLY	2.3
1	A	96	PHE	2.3
1	A	107	GLY	2.3
1	A	144	ALA	2.2
1	A	155	ILE	2.2
1	A	186	ALA	2.1
1	A	36	ARG	2.1
1	B	106	GLU	2.1
1	A	31	PRO	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.