



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:48 AM UTC

PDB ID : 1V5X / pdb_00001v5x
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

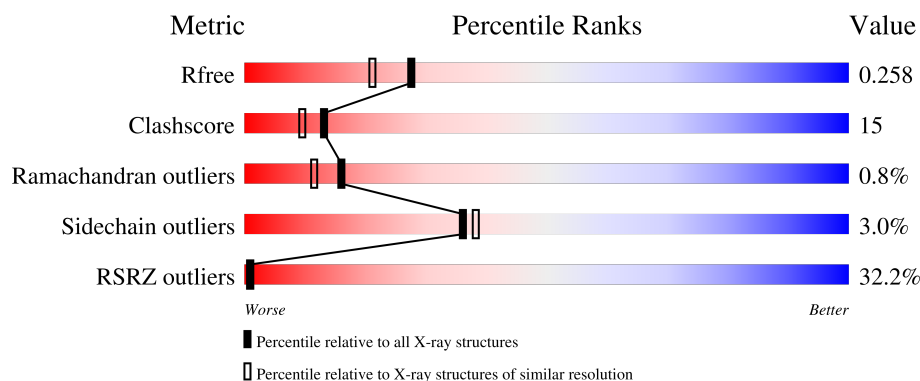
1 Overall quality at a glance

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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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 $\leq 5\%$. 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	62% 60% 34% ..
1	B	203	2% 83% 13% ..

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			

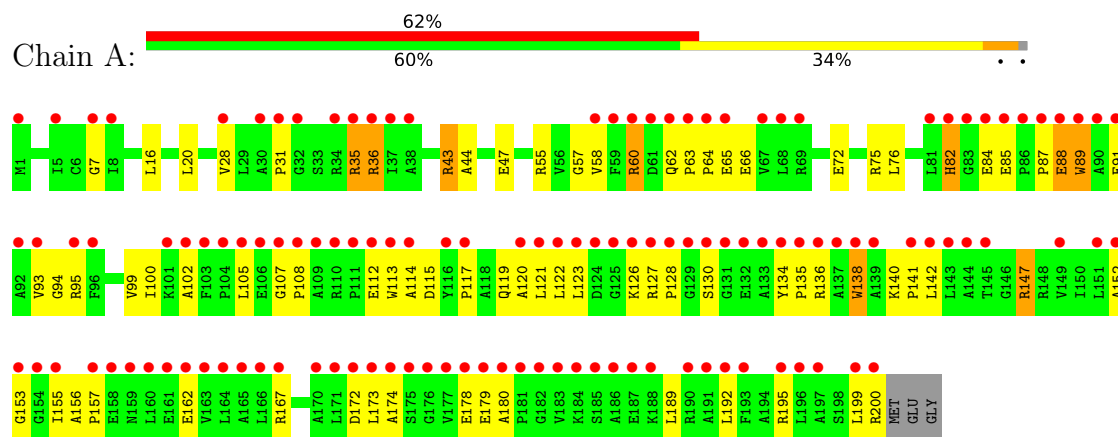
- Molecule 2 is water.

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

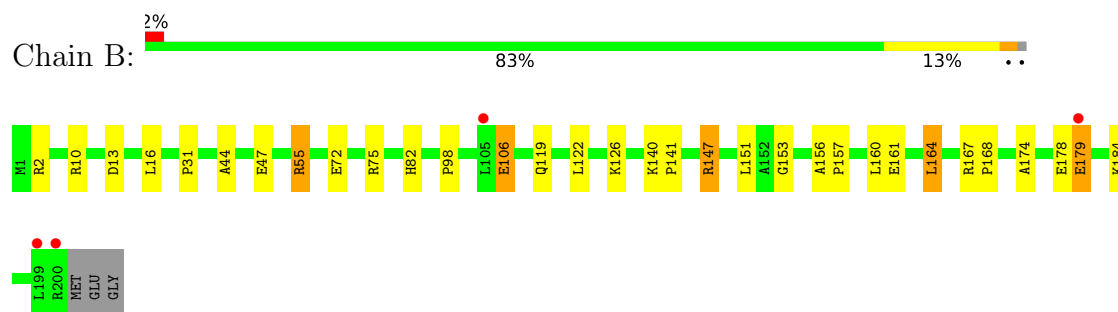
3 Residue-property plots [i](#)

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: 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, α , β , γ	144.21 Å 144.21 Å 144.21 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.00 – 2.00 35.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.00-2.00) 99.8 (35.00-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	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%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	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 ²)	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.*

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](#)

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	0.50	1/1563 (0.1%)	0.96	6/2123 (0.3%)
1	B	0.47	0/1563	0.91	3/2123 (0.1%)
All	All	0.49	1/3126 (0.0%)	0.94	9/4246 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	TRP	NE1-CE2	10.41	1.48	1.37

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	GLU	N-CA-C	-9.21	102.17	113.50
1	B	179	GLU	N-CA-C	-7.75	103.95	113.41
1	A	35	ARG	CD-NE-CZ	-5.09	117.27	124.40
1	A	36	ARG	CD-NE-CZ	-5.09	117.27	124.40
1	B	147	ARG	CD-NE-CZ	-5.08	117.28	124.40
1	A	200	ARG	CD-NE-CZ	-5.07	117.30	124.40
1	A	43	ARG	CD-NE-CZ	-5.07	117.31	124.40
1	A	60	ARG	CD-NE-CZ	-5.07	117.31	124.40
1	B	106	GLU	N-CA-C	-5.03	105.94	113.89

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:HH11	1:A:147:ARG:HB3	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:HB3	1:A:147:ARG:NH1	2.15	0.60
1:A:195:ARG:O	1:A:199:LEU:HG	2.02	0.60
1:A:113:TRP:C	1:A:115:ASP:H	2.10	0.59
1:B:55:ARG:CG	1:B:55:ARG:HH11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ARG:HH11	1:A:55:ARG:HG3	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: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
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: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:122:LEU:HD21	1:A:152:ALA:HB2	1.93	0.50
1:A:113:TRP:C	1:A:115:ASP:N	2.65	0.50
1:B:167:ARG:HH11	1:B:167:ARG:HG3	1.77	0.50
1:A:84:GLU:HA	1:A:84:GLU:OE1	2.12	0.49
1:A:114:ALA:HA	1:A:121:LEU:HD11	1.94	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:93:VAL:C	1:A:95:ARG:H	2.22	0.48
1:B:161:GLU:CD	1:B:161:GLU:H	2.22	0.47
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:122:LEU:HD23	1:B:122:LEU:C	2.41	0.46
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:HD3	1:A:127:ARG:N	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:88:GLU:H	1:A:88:GLU:CD	2.25	0.45
1:A:28:VAL:HA	1:A:58:VAL:HB	1.98	0.44
1:A:36:ARG:HH11	1:A:36:ARG:HD2	1.61	0.44
1:A:93:VAL:HG23	1:A:94:GLY:N	2.33	0.44
1:A:155:ILE:HD12	1:A:172:ASP:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:ARG:CG	1:B:55:ARG:NH1	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:57:GLY:CA	1:A:76:LEU:HD13	2.48	0.42
1:A:122:LEU:C	1:A:122:LEU:HD23	2.44	0.42
1:A:20:LEU:CD1	1:A:189:LEU:HB3	2.49	0.42
1:A:85:GLU:H	1:A:85:GLU:CD	2.25	0.42
1:A:88:GLU:CD	1:A:88:GLU:N	2.78	0.42
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:62:GLN:HB3	1:A:66:GLU:HG2	2.03	0.41
1:A:63:PRO:HA	1:A:64:PRO:HD3	1.96	0.41
1:A:105:LEU:C	1:A:107:GLY:H	2.29	0.41
1:A:140:LYS:N	1:A:141:PRO:HD2	2.36	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:HH11	1:B:147:ARG:HD2	1.63	0.40
1:A:63:PRO:HG2	1:A:66:GLU:HB2	2.03	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 ⓘ

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	198/203 (98%)	170 (86%)	26 (13%)	2 (1%)	12	8
1	B	198/203 (98%)	192 (97%)	5 (2%)	1 (0%)	24	21
All	All	396/406 (98%)	362 (91%)	31 (8%)	3 (1%)	16	11

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 ⓘ

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%)	144 (97%)	4 (3%)	39	42
1	B	148/150 (99%)	143 (97%)	5 (3%)	32	33
All	All	296/300 (99%)	287 (97%)	9 (3%)	36	38

All (9) 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	147	ARG
1	B	16	LEU
1	B	55	ARG
1	B	82	HIS
1	B	164	LEU
1	B	179	GLU

Sometimes 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 ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 ⓘ

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, 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	200/203 (98%)	2.43	125 (62%) 0 0	19, 51, 84, 89	29 (14%)
1	B	200/203 (98%)	0.33	4 (2%) 65 64	17, 30, 50, 60	4 (2%)
All	All	400/406 (98%)	1.38	129 (32%) 1 1	17, 40, 77, 89	33 (8%)

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	105	LEU	6.6
1	A	127	ARG	6.5
1	A	136	ARG	6.1
1	A	103	PHE	5.9
1	A	183	VAL	5.4
1	A	113	TRP	5.4
1	A	122	LEU	5.0
1	A	142	LEU	5.0
1	A	188	LYS	4.8
1	A	186	ALA	4.8
1	A	177	VAL	4.7
1	A	89	TRP	4.6
1	A	164	LEU	4.5
1	A	134	TYR	4.5
1	A	160	LEU	4.4
1	A	37	ILE	4.4
1	A	128	PRO	4.4
1	A	174	ALA	4.4
1	A	67	VAL	4.3
1	A	124	ASP	4.2
1	A	90	ALA	4.2
1	A	195	ARG	4.1
1	A	125	GLY	4.1
1	A	102	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	30	ALA	4.0
1	A	83	GLY	4.0
1	A	138	TRP	4.0
1	A	190	ARG	3.9
1	A	143	LEU	3.9
1	A	200	ARG	3.9
1	A	155	ILE	3.8
1	A	133	ALA	3.8
1	A	87	PRO	3.8
1	A	157	PRO	3.8
1	A	139	ALA	3.8
1	A	109	ALA	3.7
1	A	166	LEU	3.7
1	A	82	HIS	3.7
1	A	175	SER	3.6
1	A	165	ALA	3.6
1	A	84	GLU	3.6
1	A	137	ALA	3.6
1	A	110	ARG	3.6
1	A	116	TYR	3.6
1	A	192	LEU	3.5
1	A	114	ALA	3.4
1	B	200	ARG	3.4
1	A	120	ALA	3.4
1	A	179	GLU	3.4
1	A	101	LYS	3.4
1	A	95	ARG	3.4
1	A	129	GLY	3.4
1	A	131	GLY	3.4
1	A	171	LEU	3.3
1	A	185	SER	3.3
1	A	152	ALA	3.3
1	A	199	LEU	3.3
1	A	151	LEU	3.3
1	A	104	PRO	3.2
1	A	58	VAL	3.2
1	A	123	LEU	3.1
1	A	145	THR	3.1
1	A	108	PRO	3.1
1	A	69	ARG	3.1
1	A	196	LEU	3.1
1	A	141	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	181	PRO	3.1
1	A	59	PHE	3.0
1	A	184	LYS	3.0
1	A	180	ALA	3.0
1	A	68	LEU	3.0
1	A	96	PHE	3.0
1	A	35	ARG	3.0
1	A	93	VAL	3.0
1	A	86	PRO	2.9
1	A	28	VAL	2.9
1	A	135	PRO	2.9
1	A	112	GLU	2.9
1	A	191	ALA	2.9
1	A	173	LEU	2.8
1	A	111	PRO	2.8
1	A	5	ILE	2.8
1	B	105	LEU	2.8
1	A	106	GLU	2.7
1	A	38	ALA	2.7
1	A	36	ARG	2.7
1	A	197	ALA	2.7
1	A	132	GLU	2.7
1	A	161	GLU	2.7
1	A	32	GLY	2.7
1	A	170	ALA	2.6
1	A	130	SER	2.6
1	A	149	VAL	2.6
1	A	121	LEU	2.6
1	A	65	GLU	2.6
1	A	117	PRO	2.5
1	A	60	ARG	2.5
1	A	88	GLU	2.5
1	A	178	GLU	2.5
1	A	193	PHE	2.5
1	A	107	GLY	2.5
1	A	176	GLY	2.5
1	A	167	ARG	2.5
1	A	31	PRO	2.5
1	A	144	ALA	2.5
1	A	34	ARG	2.5
1	A	85	GLU	2.4
1	A	81	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	172	ASP	2.4
1	A	92	ALA	2.4
1	A	8	ILE	2.4
1	A	182	GLY	2.3
1	A	163	VAL	2.3
1	A	62	GLN	2.3
1	A	159	ASN	2.3
1	A	7	GLY	2.3
1	A	187	GLU	2.3
1	A	64	PRO	2.3
1	A	153	GLY	2.3
1	A	126	LYS	2.2
1	A	61	ASP	2.2
1	A	158	GLU	2.2
1	B	179	GLU	2.2
1	B	199	LEU	2.2
1	A	1	MET	2.1
1	A	162	GLU	2.1
1	A	154	GLY	2.1
1	A	91	GLU	2.0
1	A	63	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 oligosaccharides 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.