



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:42 PM UTC

PDB ID : 1WTJ / pdb_00001wtj
Title : Crystal Structure of delta1-piperideine-2-carboxylate reductase from *Pseudomonas syringae* pvar.tomato
Authors : Goto, M.; Muramatsu, H.; Mihara, H.; Kurihara, T.; Esaki, N.; Omi, R.; Miyahara, I.; Hirotsu, K.
Deposited on : 2004-11-24
Resolution : 1.55 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 ⓘ 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

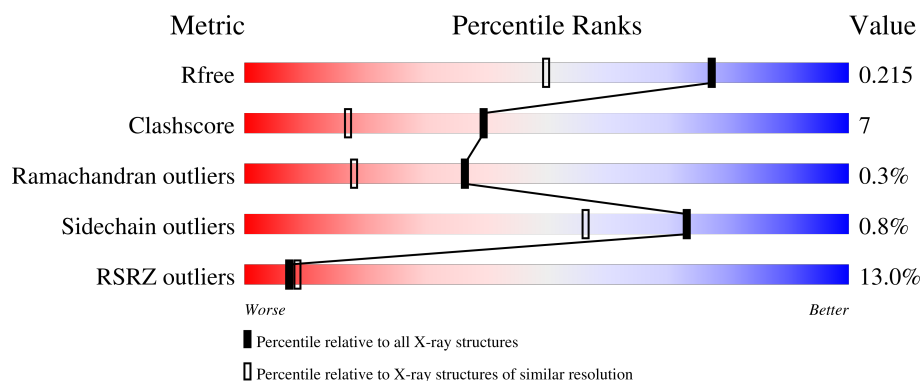
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	343	15% 82% 15% .
1	B	343	10% 82% 15% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5674 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 ureidoglycolate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	0	0
			2476	1561	452	454	9			
1	B	337	Total	C	N	O	S	0	0	0
			2514	1583	460	462	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	THR	ALA	conflict	UNP Q4U331
A	16	PRO	THR	conflict	UNP Q4U331
A	28	VAL	ALA	conflict	UNP Q4U331
A	112	ARG	LEU	conflict	UNP Q4U331
B	12	THR	ALA	conflict	UNP Q4U331
B	16	PRO	THR	conflict	UNP Q4U331
B	28	VAL	ALA	conflict	UNP Q4U331
B	112	ARG	LEU	conflict	UNP Q4U331

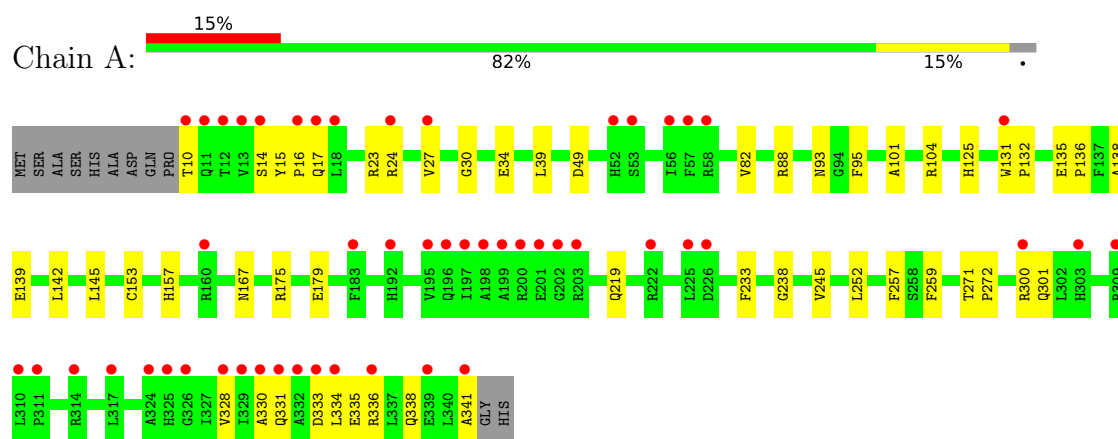
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	341	Total	O	0	0
			341	341		
2	B	343	Total	O	0	0
			343	343		

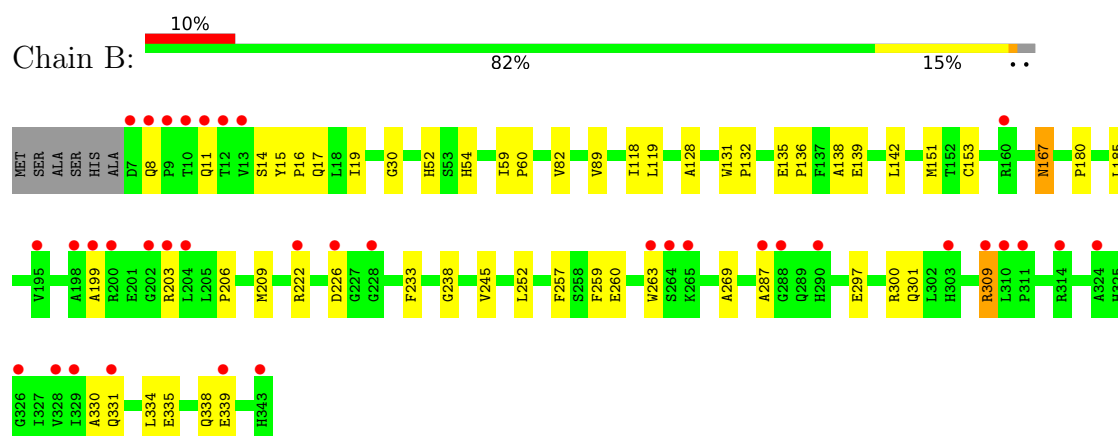
3 Residue-property plots [i](#)

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: ureidoglycolate dehydrogenase



- Molecule 1: ureidoglycolate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.44Å 68.97Å 178.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.50 – 1.55 37.50 – 1.55	Depositor EDS
% Data completeness (in resolution range)	97.7 (37.50-1.55) 97.8 (37.50-1.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.88 (at 1.55Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.194 , 0.215 0.194 , 0.215	Depositor DCC
R_{free} test set	11911 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.3	Xtriage
Anisotropy	0.658	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5674	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 4.73% 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	0.33	0/2532	0.85	4/3441 (0.1%)
1	B	0.34	0/2572	0.85	6/3496 (0.2%)
All	All	0.33	0/5104	0.85	10/6937 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	PHE	N-CA-C	-7.13	101.64	110.41
1	B	257	PHE	N-CA-C	-7.01	101.79	110.41
1	A	30	GLY	N-CA-C	6.14	123.79	115.32
1	B	30	GLY	N-CA-C	5.73	124.00	115.64
1	B	167	ASN	N-CA-C	5.47	121.91	109.81
1	B	260	GLU	N-CA-C	5.27	118.53	112.57
1	B	52	HIS	N-CA-C	5.26	117.69	111.33
1	A	82	VAL	N-CA-C	5.23	117.92	112.90
1	B	82	VAL	N-CA-C	5.11	117.89	112.83
1	A	219	GLN	N-CA-C	-5.00	107.21	113.72

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	2476	0	2452	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2514	0	2481	37	0
2	A	341	0	0	1	0
2	B	343	0	0	2	0
All	All	5674	0	4933	66	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 7.

All (66) 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:24:ARG:HD2	1:A:341:ALA:HB1	1.41	0.99
1:B:206:PRO:HG2	1:B:209:MET:SD	2.36	0.65
1:A:101:ALA:HA	1:A:104:ARG:NH1	2.13	0.63
1:B:222:ARG:HD2	1:B:226:ASP:OD2	2.00	0.61
1:A:14:SER:OG	1:A:17:GLN:HG3	2.02	0.59
1:B:14:SER:OG	1:B:17:GLN:HG3	2.03	0.58
1:A:101:ALA:HA	1:A:104:ARG:HH12	1.69	0.58
1:B:263:TRP:HB2	1:B:269:ALA:HB3	1.85	0.57
1:B:11:GLN:NE2	1:B:334:LEU:HD23	2.20	0.57
1:B:297:GLU:HA	1:B:300:ARG:NH1	2.20	0.56
1:A:252:LEU:C	1:A:252:LEU:HD23	2.30	0.56
1:A:331:GLN:O	1:A:335:GLU:HG3	2.06	0.56
1:A:233:PHE:O	1:A:238:GLY:HA3	2.08	0.54
1:B:54:HIS:HD2	2:B:501:HOH:O	1.91	0.53
1:A:334:LEU:O	1:A:338:GLN:HG3	2.08	0.53
1:A:125:HIS:HD2	1:A:271:THR:OG1	1.92	0.53
1:A:301:GLN:HG3	1:B:259:PHE:CE1	2.45	0.52
1:A:135:GLU:HB3	1:A:136:PRO:HD3	1.90	0.52
1:A:259:PHE:CE1	1:B:301:GLN:HG3	2.45	0.51
1:A:49:ASP:OD2	1:A:131:TRP:HH2	1.93	0.51
1:B:11:GLN:OE1	1:B:331:GLN:HG3	2.11	0.51
1:B:167:ASN:HB3	1:B:245:VAL:HG11	1.93	0.51
1:B:331:GLN:O	1:B:335:GLU:HG3	2.11	0.51
1:B:334:LEU:O	1:B:338:GLN:HG3	2.11	0.50
1:B:8:GLN:HG2	1:B:330:ALA:HB1	1.92	0.50
1:B:135:GLU:HB3	1:B:136:PRO:HD3	1.93	0.50
1:A:300:ARG:HH21	1:A:301:GLN:NE2	2.10	0.49
1:B:199:ALA:HB1	1:B:222:ARG:NH1	2.26	0.49
1:A:93:ASN:OD1	1:A:125:HIS:HE1	1.96	0.49
1:A:49:ASP:OD2	1:A:131:TRP:CH2	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:TYR:N	1:B:16:PRO:HD2	2.28	0.47
1:A:15:TYR:N	1:A:16:PRO:HD2	2.29	0.47
1:A:131:TRP:CG	1:A:132:PRO:HD3	2.50	0.46
1:B:300:ARG:CZ	1:B:300:ARG:HB2	2.46	0.46
1:A:167:ASN:HB3	1:A:245:VAL:HG11	1.98	0.46
1:A:34:GLU:H	1:A:34:GLU:CD	2.24	0.46
1:A:272:PRO:HG3	2:A:533:HOH:O	2.15	0.46
1:B:59:ILE:HB	1:B:60:PRO:HD3	1.98	0.46
1:B:233:PHE:O	1:B:238:GLY:HA3	2.17	0.45
1:A:10:THR:HG22	1:A:330:ALA:HA	1.98	0.45
1:B:309:ARG:HD3	1:B:309:ARG:H	1.81	0.45
1:A:333:ASP:O	1:A:336:ARG:HG2	2.17	0.44
1:B:89:VAL:HG23	1:B:119:LEU:HD11	1.98	0.44
1:B:297:GLU:HA	1:B:300:ARG:HH12	1.81	0.44
1:A:88:ARG:NH2	1:B:287:ALA:HB2	2.33	0.44
1:B:252:LEU:HD23	1:B:252:LEU:C	2.43	0.43
1:A:300:ARG:HH21	1:A:301:GLN:HE21	1.66	0.43
1:B:185:LEU:HD12	1:B:185:LEU:O	2.19	0.42
1:B:128:ALA:HB1	2:B:353:HOH:O	2.18	0.42
1:B:14:SER:HG	1:B:17:GLN:HG3	1.83	0.42
1:A:175:ARG:HD3	1:A:179:GLU:O	2.20	0.42
1:B:15:TYR:CZ	1:B:19:ILE:HD11	2.55	0.42
1:A:138:ALA:HA	1:A:142:LEU:O	2.19	0.42
1:B:131:TRP:N	1:B:132:PRO:CD	2.83	0.42
1:B:309:ARG:NE	1:B:309:ARG:O	2.53	0.41
1:A:157:HIS:O	1:B:309:ARG:NH1	2.53	0.41
1:A:135:GLU:O	1:A:139:GLU:HG3	2.20	0.41
1:B:135:GLU:O	1:B:139:GLU:HG3	2.19	0.41
1:B:138:ALA:HB1	1:B:180:PRO:HB3	2.01	0.41
1:A:39:LEU:HD11	1:A:95:PHE:CD2	2.55	0.41
1:A:328:VAL:HG13	1:A:328:VAL:O	2.20	0.41
1:B:335:GLU:O	1:B:339:GLU:HG3	2.20	0.41
1:A:131:TRP:N	1:A:132:PRO:CD	2.84	0.40
1:B:138:ALA:HA	1:B:142:LEU:O	2.21	0.40
1:B:203:ARG:HH11	1:B:203:ARG:HG3	1.85	0.40
1:A:23:ARG:O	1:A:27:VAL:HG23	2.22	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	330/343 (96%)	323 (98%)	6 (2%)	1 (0%)	36	18
1	B	335/343 (98%)	326 (97%)	8 (2%)	1 (0%)	36	18
All	All	665/686 (97%)	649 (98%)	14 (2%)	2 (0%)	36	18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	153	CYS
1	B	153	CYS

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	250/258 (97%)	249 (100%)	1 (0%)	84	73
1	B	254/258 (98%)	251 (99%)	3 (1%)	63	40
All	All	504/516 (98%)	500 (99%)	4 (1%)	73	56

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LEU
1	B	118	ILE
1	B	151	MET
1	B	309	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	HIS
1	A	290	HIS
1	A	301	GLN
1	B	8	GLN
1	B	11	GLN
1	B	54	HIS
1	B	192	HIS
1	B	276	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 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	332/343 (96%)	0.68	51 (15%) 5 6	12, 19, 37, 51	0
1	B	337/343 (98%)	0.54	36 (10%) 11 13	12, 18, 34, 46	0
All	All	669/686 (97%)	0.61	87 (13%) 7 9	12, 19, 36, 51	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	ILE	6.0
1	B	309	ARG	4.8
1	A	330	ALA	4.7
1	A	334	LEU	4.4
1	A	57	PHE	4.2
1	A	310	LEU	4.1
1	B	314	ARG	3.8
1	A	336	ARG	3.8
1	B	203	ARG	3.8
1	A	10	THR	3.8
1	B	310	LEU	3.7
1	B	8	GLN	3.7
1	A	332	ALA	3.7
1	A	324	ALA	3.5
1	B	13	VAL	3.5
1	A	201	GLU	3.4
1	B	12	THR	3.3
1	A	199	ALA	3.3
1	A	12	THR	3.3
1	A	13	VAL	3.3
1	A	52	HIS	3.3
1	A	328	VAL	3.2
1	A	200	ARG	3.2
1	B	287	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	288	GLY	3.2
1	A	197	ILE	3.1
1	B	263	TRP	3.1
1	B	222	ARG	3.0
1	A	341	ALA	3.0
1	B	202	GLY	3.0
1	B	11	GLN	3.0
1	A	11	GLN	2.9
1	A	24	ARG	2.9
1	B	9	PRO	2.9
1	A	317	LEU	2.9
1	B	228	GLY	2.9
1	A	16	PRO	2.8
1	A	222	ARG	2.8
1	B	160	ARG	2.6
1	A	196	GLN	2.6
1	A	198	ALA	2.6
1	A	331	GLN	2.6
1	A	325	HIS	2.6
1	B	226	ASP	2.6
1	B	329	ILE	2.5
1	A	27	VAL	2.5
1	B	343	HIS	2.5
1	A	203	ARG	2.5
1	A	303	HIS	2.5
1	A	339	GLU	2.5
1	A	311	PRO	2.5
1	A	195	VAL	2.5
1	A	56	ILE	2.4
1	A	225	LEU	2.4
1	A	226	ASP	2.4
1	A	18	LEU	2.4
1	A	131	TRP	2.4
1	A	160	ARG	2.4
1	B	311	PRO	2.4
1	B	331	GLN	2.4
1	B	198	ALA	2.3
1	A	300	ARG	2.3
1	B	303	HIS	2.3
1	B	10	THR	2.3
1	A	314	ARG	2.3
1	B	199	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	324	ALA	2.3
1	B	328	VAL	2.3
1	A	202	GLY	2.3
1	A	192	HIS	2.2
1	B	290	HIS	2.2
1	A	58	ARG	2.2
1	A	183	PHE	2.2
1	B	265	LYS	2.2
1	B	264	SER	2.2
1	B	200	ARG	2.2
1	A	326	GLY	2.1
1	A	53	SER	2.1
1	A	14	SER	2.1
1	B	326	GLY	2.1
1	B	7	ASP	2.1
1	A	333	ASP	2.1
1	A	309	ARG	2.0
1	A	17	GLN	2.0
1	B	339	GLU	2.0
1	B	195	VAL	2.0
1	B	204	LEU	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.