



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2022 – 08:11 PM JST

PDB ID : 7W0I  
Title : Crystal structure of threonine aldolase from *Mycobacterium vanbaalenii*  
Authors : Wu, B.  
Deposited on : 2021-11-18  
Resolution : 1.75 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

<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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
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)	:	2.31.2

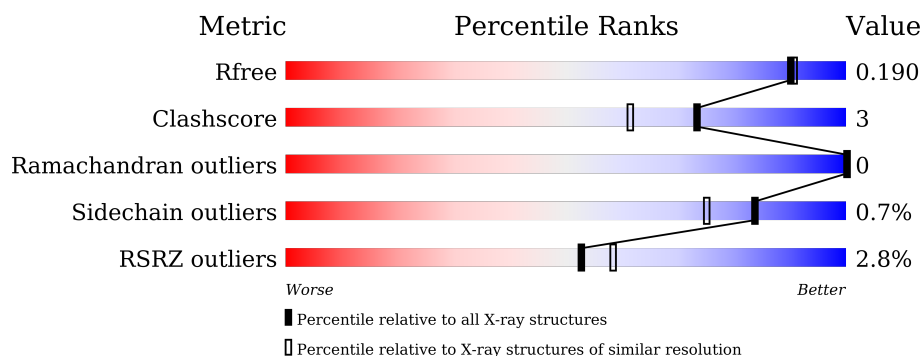
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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  $\geq 3$ , 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	<div> <div>3%</div> <div>92%</div> <div>5%</div> <div>.</div> </div>
1	B	343	<div> <div>3%</div> <div>92%</div> <div>5%</div> <div>.</div> </div>
1	C	343	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	D	343	<div> <div>3%</div> <div>90%</div> <div>7%</div> <div>.</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11304 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 L-threonine aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2504	1567	466	467	4			
1	B	334	Total	C	N	O	S	0	0	0
			2504	1567	466	467	4			
1	C	334	Total	C	N	O	S	0	0	0
			2504	1567	466	467	4			
1	D	334	Total	C	N	O	S	0	0	0
			2504	1567	466	467	4			

- Molecule 2 is water.

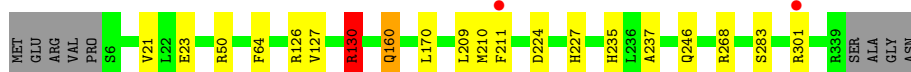
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	362	Total	O	0	0
			362	362		
2	B	338	Total	O	0	0
			338	338		
2	C	323	Total	O	0	0
			323	323		
2	D	265	Total	O	0	0
			265	265		

### 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: L-threonine aldolase

Chain A: 

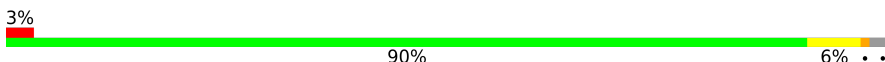


- Molecule 1: L-threonine aldolase

Chain B: 




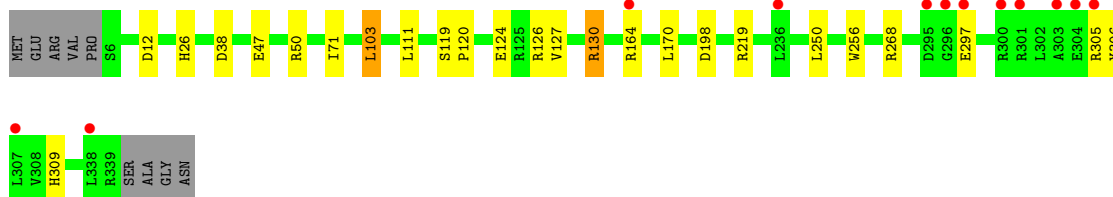
- Molecule 1: L-threonine aldolase

Chain C: 



- Molecule 1: L-threonine aldolase

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.00Å 102.43Å 150.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.48 – 1.75 48.48 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.48-1.75) 99.8 (48.48-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.53 (at 1.75Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.171 , 0.190 0.171 , 0.190	Depositor DCC
$R_{free}$ test set	1998 reflections (1.34%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.0	Xtriage
Anisotropy	0.499	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 46.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	11304	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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

### 5.1 Standard geometry

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	0.40	0/2559	0.60	3/3487 (0.1%)
1	B	0.41	0/2559	0.61	2/3487 (0.1%)
1	C	0.42	1/2559 (0.0%)	0.72	8/3487 (0.2%)
1	D	0.38	0/2559	0.69	10/3487 (0.3%)
All	All	0.40	1/10236 (0.0%)	0.66	23/13948 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	305	ARG	CB-CG	5.47	1.67	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	305	ARG	NE-CZ-NH2	12.11	126.36	120.30
1	D	130	ARG	NE-CZ-NH2	-9.15	115.73	120.30
1	D	130	ARG	NE-CZ-NH1	8.95	124.77	120.30
1	C	301	ARG	NE-CZ-NH2	-8.68	115.96	120.30
1	D	268	ARG	NE-CZ-NH2	-8.11	116.25	120.30
1	C	301	ARG	NE-CZ-NH1	8.07	124.34	120.30
1	D	268	ARG	CG-CD-NE	-7.96	95.09	111.80
1	C	130	ARG	NE-CZ-NH2	-7.91	116.35	120.30
1	D	268	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	C	130	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	D	103	LEU	CB-CG-CD2	7.04	122.97	111.00
1	C	339	ARG	NE-CZ-NH2	-6.71	116.94	120.30
1	D	297	GLU	CA-CB-CG	6.62	127.98	113.40
1	A	130	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	D	38	ASP	CB-CG-OD2	-5.93	112.96	118.30
1	C	297	GLU	CA-CB-CG	5.88	126.34	113.40
1	C	301	ARG	CB-CA-C	-5.85	98.70	110.40
1	A	160	GLN	CA-CB-CG	5.82	126.19	113.40
1	A	130	ARG	NE-CZ-NH1	5.22	122.91	120.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	103	LEU	CB-CA-C	-5.21	100.29	110.20
1	D	38	ASP	CB-CG-OD1	5.19	122.97	118.30
1	B	338	LEU	CB-CG-CD2	5.18	119.81	111.00
1	B	280	GLN	CA-CB-CG	5.08	124.58	113.40

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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	2504	0	2456	18	0
1	B	2504	0	2456	19	0
1	C	2504	0	2456	16	0
1	D	2504	0	2456	16	0
2	A	362	0	0	9	1
2	B	338	0	0	8	2
2	C	323	0	0	9	2
2	D	265	0	0	6	1
All	All	11304	0	9824	65	3

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 3.

All (65) 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:268:ARG:NH2	2:C:403:HOH:O	2.01	0.91
1:C:306:TYR:OH	2:C:401:HOH:O	1.95	0.84
1:B:339:ARG:NH2	2:B:401:HOH:O	1.81	0.80
1:C:313:PRO:O	2:C:402:HOH:O	1.99	0.79
1:B:301:ARG:NH1	2:B:402:HOH:O	2.17	0.77
1:D:124:GLU:OE1	1:D:164:ARG:NH2	2.18	0.76
1:B:296:GLY:O	1:B:300:ARG:HG3	1.86	0.75
1:D:130:ARG:HD3	2:D:421:HOH:O	1.86	0.74

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:ARG:NH1	2:B:403:HOH:O	2.21	0.73
1:C:6:SER:N	2:C:405:HOH:O	2.22	0.71
1:C:224:ASP:O	2:C:404:HOH:O	2.09	0.70
1:A:283:SER:OG	2:A:401:HOH:O	2.09	0.69
1:D:47:GLU:OE1	2:D:401:HOH:O	2.10	0.68
1:B:302:LEU:HD22	1:B:338:LEU:HD21	1.77	0.66
1:A:23:GLU:OE2	2:A:402:HOH:O	2.13	0.66
1:A:130:ARG:HD3	2:A:422:HOH:O	1.97	0.64
1:A:160:GLN:NE2	2:A:404:HOH:O	2.29	0.64
1:D:126:ARG:HD3	2:D:432:HOH:O	1.97	0.62
1:C:130:ARG:HD3	2:C:420:HOH:O	1.99	0.62
1:C:295:ASP:OD2	1:C:297:GLU:HB2	2.01	0.60
1:B:64:PHE:H	1:B:235:HIS:HD2	1.49	0.60
1:B:79:ARG:NH1	2:B:405:HOH:O	2.37	0.58
1:A:64:PHE:H	1:A:235:HIS:HD2	1.52	0.58
1:D:127:VAL:HG21	1:D:170:LEU:HD11	1.86	0.57
1:C:296:GLY:O	1:C:300:ARG:HG3	2.05	0.56
1:C:314:GLY:C	1:C:315:GLN:HG2	2.26	0.56
1:B:130:ARG:HD3	2:B:417:HOH:O	2.07	0.55
1:B:50:ARG:NH2	2:B:408:HOH:O	2.42	0.53
1:B:64:PHE:H	1:B:235:HIS:CD2	2.27	0.53
1:D:309:HIS:HE1	2:D:477:HOH:O	1.92	0.53
1:A:211:PHE:CE2	1:B:211:PHE:HE2	2.28	0.51
1:C:111:LEU:HD23	1:C:126:ARG:HD2	1.93	0.50
1:B:50:ARG:HD3	1:B:58:ALA:O	2.10	0.50
1:A:21:VAL:HG13	1:A:246:GLN:HG2	1.94	0.50
1:C:22:LEU:HD13	1:D:26:HIS:CE1	2.48	0.49
1:A:127:VAL:HG21	1:A:170:LEU:HD21	1.93	0.48
1:A:210:MET:HG2	1:B:242:PHE:CG	2.48	0.48
1:D:71:ILE:HG22	1:D:103:LEU:HD22	1.94	0.48
1:D:198:ASP:OD1	1:D:219:ARG:HD2	2.14	0.48
1:A:126:ARG:HD3	2:A:556:HOH:O	2.14	0.47
1:A:235:HIS:HE1	1:B:233:ILE:O	1.97	0.47
1:D:250:LEU:HG	1:D:256:TRP:HB3	1.96	0.47
1:B:268:ARG:O	1:B:272:THR:HG23	2.15	0.47
1:C:309:HIS:HE1	2:C:472:HOH:O	1.97	0.47
1:A:268:ARG:HE	1:A:268:ARG:HB3	1.41	0.46
1:A:160:GLN:CD	2:A:404:HOH:O	2.54	0.46
1:A:227:HIS:HD2	2:A:530:HOH:O	1.99	0.45
1:C:64:PHE:CZ	1:C:237:ALA:HB2	2.52	0.45
1:D:50:ARG:NH2	2:D:413:HOH:O	2.49	0.45

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:PHE:CZ	1:A:237:ALA:HB2	2.51	0.45
1:D:306:TYR:OH	2:D:402:HOH:O	2.21	0.44
1:D:12:ASP:OD1	1:D:12:ASP:N	2.50	0.43
1:D:111:LEU:HD23	1:D:126:ARG:HD2	2.00	0.43
1:A:209:LEU:HD22	1:A:246:GLN:HB3	2.00	0.43
1:A:50:ARG:NH2	2:A:414:HOH:O	2.52	0.42
1:B:272:THR:HG22	2:B:710:HOH:O	2.19	0.42
1:C:268:ARG:NH1	2:C:408:HOH:O	2.43	0.42
1:B:64:PHE:CZ	1:B:237:ALA:HB2	2.55	0.42
1:D:164:ARG:HH11	1:D:164:ARG:HG3	1.85	0.42
1:C:127:VAL:HG21	1:C:170:LEU:HD21	2.02	0.41
1:B:280:GLN:HB2	2:B:447:HOH:O	2.20	0.41
1:D:120:PRO:HB3	1:D:164:ARG:HG3	2.02	0.41
1:C:50:ARG:NH2	2:C:414:HOH:O	2.53	0.41
1:B:300:ARG:HH11	1:B:300:ARG:HD2	1.49	0.40
1:A:224:ASP:O	2:A:403:HOH:O	2.22	0.40

All (3) 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)
2:A:670:HOH:O	2:D:658:HOH:O[3_545]	1.93	0.27
2:B:654:HOH:O	2:C:680:HOH:O[3_445]	1.94	0.26
2:B:675:HOH:O	2:C:614:HOH:O[3_445]	2.07	0.13

## 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	332/343 (97%)	325 (98%)	7 (2%)	0	100	100
1	B	332/343 (97%)	325 (98%)	7 (2%)	0	100	100

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	332/343 (97%)	324 (98%)	8 (2%)	0	100	100
1	D	332/343 (97%)	325 (98%)	7 (2%)	0	100	100
All	All	1328/1372 (97%)	1299 (98%)	29 (2%)	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	253/260 (97%)	251 (99%)	2 (1%)	81	72
1	B	253/260 (97%)	252 (100%)	1 (0%)	91	87
1	C	253/260 (97%)	251 (99%)	2 (1%)	81	72
1	D	253/260 (97%)	251 (99%)	2 (1%)	81	72
All	All	1012/1040 (97%)	1005 (99%)	7 (1%)	84	75

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	ARG
1	A	301	ARG
1	B	130	ARG
1	C	130	ARG
1	C	300	ARG
1	D	119	SER
1	D	305	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	GLN
1	A	227	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	235	HIS
1	B	235	HIS
1	B	280	GLN
1	C	309	HIS

### 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 monosaccharides 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, 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	334/343 (97%)	-0.22	2 (0%) 89 92	14, 21, 38, 61	0
1	B	334/343 (97%)	-0.06	12 (3%) 42 49	15, 21, 40, 74	0
1	C	334/343 (97%)	-0.05	12 (3%) 42 49	16, 26, 50, 84	0
1	D	334/343 (97%)	0.15	12 (3%) 42 49	17, 32, 48, 79	0
All	All	1336/1372 (97%)	-0.04	38 (2%) 53 58	14, 25, 45, 84	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	300	ARG	4.9
1	C	339	ARG	4.8
1	B	300	ARG	4.8
1	B	295	ASP	4.7
1	D	301	ARG	4.7
1	B	296	GLY	4.6
1	B	297	GLU	4.4
1	C	296	GLY	4.3
1	C	295	ASP	4.2
1	C	304	GLU	4.2
1	D	297	GLU	4.1
1	C	301	ARG	4.1
1	B	339	ARG	3.9
1	D	300	ARG	3.8
1	C	297	GLU	3.8
1	D	303	ALA	3.7
1	C	303	ALA	3.6
1	D	296	GLY	3.6
1	D	307	LEU	3.5
1	D	305	ARG	3.2
1	B	338	LEU	3.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	305	ARG	3.1
1	B	301	ARG	3.1
1	D	295	ASP	3.1
1	B	195	GLY	3.0
1	C	337	ALA	2.9
1	D	338	LEU	2.7
1	C	336	ALA	2.7
1	B	314	GLY	2.7
1	C	278	GLY	2.7
1	B	303	ALA	2.6
1	A	211	PHE	2.5
1	D	304	GLU	2.4
1	D	236	LEU	2.3
1	A	301	ARG	2.3
1	B	304	GLU	2.3
1	B	298	ALA	2.2
1	D	164	ARG	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 monosaccharides 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.