



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 14, 2022 – 04:08 PM JST

PDB ID : 7XYW  
Title : Crystal structure of ZYG11B bound to AFLH degron  
Authors : Dong, C.; Yan, X.; Li, Y.  
Deposited on : 2022-06-02  
Resolution : 2.50 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)

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<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>.

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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)	:	1.13
EDS	:	2.31.3
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)	:	2.31.3

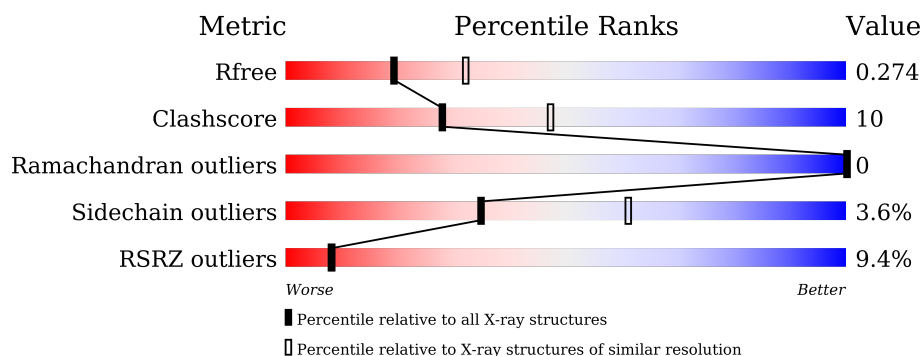
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	250	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>22%</div> <div>..</div> </div> </div>
1	B	250	<div> <div>13%</div> <div> <div></div> <div>72%</div> <div>24%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3920 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 Protein zyg-11 homolog B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	245	Total	C	N	O	S	0	0	0
			1961	1256	333	362	10			
1	A	244	Total	C	N	O	S	0	0	0
			1955	1253	332	361	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	479	ALA	-	expression tag	UNP Q9C0D3
B	480	PHE	-	expression tag	UNP Q9C0D3
B	481	LEU	-	expression tag	UNP Q9C0D3
B	482	HIS	-	expression tag	UNP Q9C0D3
B	483	VAL	-	expression tag	UNP Q9C0D3
B	484	GLY	-	expression tag	UNP Q9C0D3
A	479	ALA	-	expression tag	UNP Q9C0D3
A	480	PHE	-	expression tag	UNP Q9C0D3
A	481	LEU	-	expression tag	UNP Q9C0D3
A	482	HIS	-	expression tag	UNP Q9C0D3
A	483	VAL	-	expression tag	UNP Q9C0D3
A	484	GLY	-	expression tag	UNP Q9C0D3

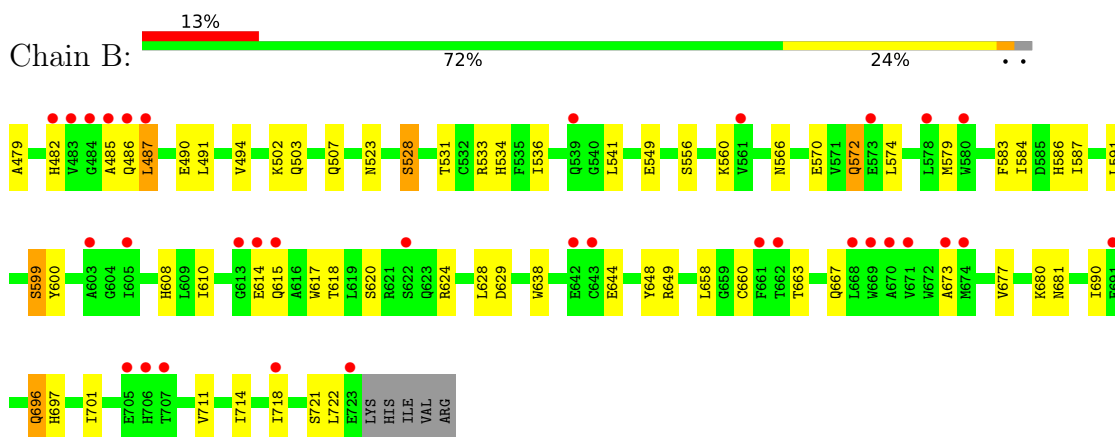
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		

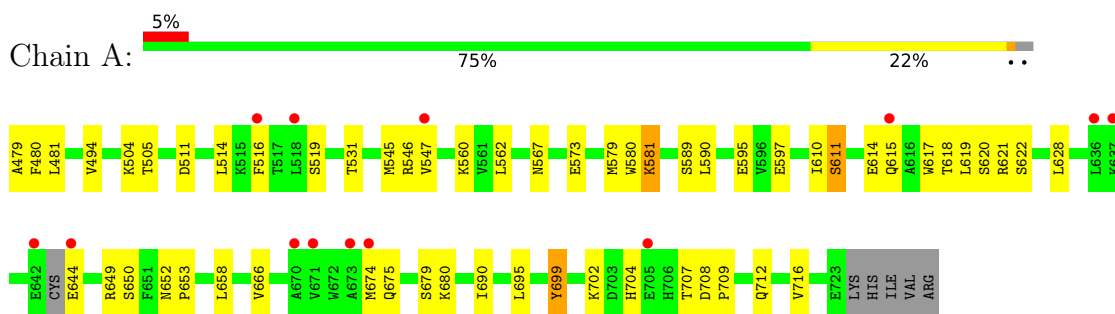
### 3 Residue-property plots [i](#)

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: Protein zyg-11 homolog B



#### • Molecule 1: Protein zyg-11 homolog B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.88Å 97.27Å 122.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.49 – 2.50 20.49 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.0 (20.49-2.50) 92.0 (20.49-2.50)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 2.50Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, $R_{free}$	0.226 , 0.275 0.225 , 0.274	Depositor DCC
$R_{free}$ test set	1039 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	67.3	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 64.9	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	0.95	EDS
Total number of atoms	3920	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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	0.54	1/2002 (0.0%)	0.66	0/2722
1	B	0.50	0/2009	0.63	2/2733 (0.1%)
All	All	0.52	1/4011 (0.0%)	0.64	2/5455 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	699	TYR	CD1-CE1	5.57	1.47	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	696	GLN	CA-CB-CG	-6.23	99.69	113.40
1	B	541	LEU	CA-CB-CG	5.74	128.50	115.30

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	1955	0	1925	38	0
1	B	1961	0	1931	45	0
2	A	4	0	0	1	0
All	All	3920	0	3856	77	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 10.

All (77) 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:649:ARG:NH1	2:A:801:HOH:O	2.04	0.89
1:A:614:GLU:HB2	1:A:621:ARG:HH22	1.53	0.72
1:B:503:GLN:HE21	1:B:507:GLN:HE22	1.39	0.69
1:A:704:HIS:O	1:A:707:THR:HG23	1.97	0.64
1:A:581:LYS:H	1:A:581:LYS:HD2	1.64	0.62
1:B:503:GLN:HE21	1:B:507:GLN:NE2	1.97	0.62
1:A:614:GLU:H	1:A:614:GLU:CD	2.02	0.61
1:B:697:HIS:O	1:B:701:ILE:HG12	2.02	0.60
1:A:579:MET:HE1	1:A:617:TRP:HA	1.84	0.60
1:B:482:HIS:ND1	1:B:485:ALA:HB2	2.15	0.60
1:B:644:GLU:HG2	1:B:680:LYS:HB3	1.84	0.60
1:A:614:GLU:HB2	1:A:621:ARG:NH2	2.17	0.58
1:A:675:GLN:O	1:A:679:SER:OG	2.17	0.56
1:A:579:MET:HE2	1:A:618:THR:H	1.71	0.55
1:A:562:LEU:HD11	1:A:590:LEU:HD13	1.90	0.54
1:A:699:TYR:O	1:A:702:LYS:HB3	2.08	0.54
1:A:595:GLU:HG3	1:A:597:GLU:OE2	2.07	0.53
1:B:617:TRP:CG	1:B:624:ARG:HD2	2.43	0.53
1:B:617:TRP:NE1	1:B:620:SER:O	2.42	0.52
1:B:486:GLN:HG3	1:B:490:GLU:OE2	2.10	0.52
1:B:523:ASN:ND2	1:A:480:PHE:HB2	2.25	0.52
1:B:644:GLU:OE1	1:B:681:ASN:HB2	2.10	0.51
1:B:523:ASN:HD22	1:A:480:PHE:HB2	1.75	0.51
1:B:600:TYR:HA	1:B:638:TRP:CH2	2.45	0.51
1:B:696:GLN:NE2	1:B:696:GLN:HA	2.24	0.51
1:B:523:ASN:HD21	1:A:481:LEU:H	1.59	0.50
1:B:721:SER:O	1:B:722:LEU:HD12	2.12	0.50
1:B:523:ASN:ND2	1:A:481:LEU:H	2.10	0.50
1:B:690:ILE:HD11	1:B:722:LEU:HD23	1.94	0.49
1:A:690:ILE:HG12	1:A:695:LEU:HD11	1.95	0.49
1:B:570:GLU:HG2	1:B:608:HIS:CE1	2.47	0.49
1:B:479:ALA:N	1:A:567:ASN:OD1	2.46	0.49
1:B:491:LEU:HB3	1:B:534:HIS:NE2	2.27	0.49
1:B:528:SER:O	1:B:528:SER:OG	2.26	0.49
1:B:714:ILE:O	1:B:718:ILE:HG13	2.12	0.49
1:B:487:LEU:HA	1:B:490:GLU:CD	2.33	0.48
1:B:503:GLN:NE2	1:B:507:GLN:OE1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610:ILE:HD13	1:A:628:LEU:HB3	1.95	0.48
1:B:482:HIS:CE1	1:B:485:ALA:HB2	2.49	0.48
1:B:494:VAL:HG21	1:B:531:THR:HG23	1.96	0.47
1:A:644:GLU:N	1:A:680:LYS:HZ1	2.13	0.47
1:A:658:LEU:HD21	1:A:674:MET:HE3	1.97	0.47
1:B:583:PHE:O	1:B:587:ILE:HG12	2.14	0.47
1:A:516:PHE:CD1	1:A:516:PHE:N	2.83	0.46
1:B:673:ALA:O	1:B:677:VAL:HG22	2.16	0.46
1:A:504:LYS:HG3	1:A:514:LEU:HB2	1.98	0.46
1:A:581:LYS:HE3	1:A:619:LEU:HD21	1.97	0.46
1:B:721:SER:C	1:B:722:LEU:HD12	2.37	0.46
1:A:494:VAL:HG21	1:A:531:THR:HG23	1.98	0.45
1:B:536:ILE:HD12	1:B:574:LEU:HD22	1.98	0.45
1:B:660:CYS:HB3	1:B:663:THR:OG1	2.17	0.45
1:A:712:GLN:O	1:A:716:VAL:HG13	2.15	0.45
1:B:591:LEU:O	1:B:599:SER:HB2	2.15	0.45
1:A:611:SER:HB2	1:A:666:VAL:HG11	1.98	0.44
1:B:502:LYS:HG3	1:B:503:GLN:N	2.33	0.44
1:B:491:LEU:HD13	1:B:534:HIS:CG	2.53	0.44
1:B:533:ARG:HD2	1:B:574:LEU:HD21	2.00	0.44
1:A:579:MET:CE	1:A:617:TRP:HA	2.48	0.44
1:B:610:ILE:HD12	1:B:628:LEU:HD12	2.00	0.44
1:B:648:TYR:CE1	1:A:479:ALA:HB2	2.53	0.43
1:B:721:SER:O	1:B:721:SER:OG	2.34	0.43
1:B:549:GLU:OE2	1:B:586:HIS:HE1	2.01	0.43
1:A:504:LYS:HD2	1:A:511:ASP:OD2	2.18	0.43
1:B:572:GLN:CD	1:B:572:GLN:H	2.22	0.42
1:B:718:ILE:O	1:B:722:LEU:HD13	2.20	0.42
1:A:614:GLU:HG2	1:A:615:GLN:H	1.84	0.42
1:A:650:SER:OG	1:A:652:ASN:OD1	2.38	0.42
1:B:583:PHE:HD2	1:B:584:ILE:HD13	1.86	0.41
1:B:556:SER:O	1:B:560:LYS:HG3	2.20	0.41
1:A:505:THR:OG1	1:A:547:VAL:HG12	2.21	0.41
1:A:545:MET:HG3	1:A:580:TRP:CZ3	2.55	0.41
1:A:708:ASP:HA	1:A:709:PRO:HD3	1.92	0.41
1:B:579:MET:HG3	1:B:618:THR:HG23	2.02	0.41
1:A:519:SER:OG	1:A:560:LYS:HE2	2.21	0.41
1:A:573:GLU:H	1:A:573:GLU:CD	2.22	0.41
1:A:652:ASN:N	1:A:653:PRO:HD2	2.37	0.40
1:B:667:GLN:HB3	1:B:711:VAL:HG11	2.03	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	240/250 (96%)	234 (98%)	6 (2%)	0	100	100
1	B	243/250 (97%)	232 (96%)	11 (4%)	0	100	100
All	All	483/500 (97%)	466 (96%)	17 (4%)	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	219/226 (97%)	213 (97%)	6 (3%)	44	71
1	B	220/226 (97%)	210 (96%)	10 (4%)	27	51
All	All	439/452 (97%)	423 (96%)	16 (4%)	35	61

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	487	LEU
1	B	528	SER
1	B	566	ASN
1	B	572	GLN
1	B	599	SER
1	B	614	GLU
1	B	615	GLN
1	B	629	ASP

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Mol	Chain	Res	Type
1	B	649	ARG
1	B	658	LEU
1	A	546	ARG
1	A	581	LYS
1	A	589	SER
1	A	611	SER
1	A	620	SER
1	A	622	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	503	GLN
1	B	507	GLN
1	B	523	ASN
1	B	538	ASN
1	B	586	HIS
1	B	696	GLN
1	A	615	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 monosaccharides 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	244/250 (97%)	0.36	13 (5%) 26 28	52, 80, 102, 113	0
1	B	245/250 (98%)	0.68	33 (13%) 3 2	55, 79, 105, 132	0
All	All	489/500 (97%)	0.52	46 (9%) 8 8	52, 79, 104, 132	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	705	GLU	5.2
1	A	615	GLN	4.9
1	B	487	LEU	4.6
1	B	706	HIS	4.4
1	B	486	GLN	4.3
1	B	484	GLY	3.9
1	B	482	HIS	3.8
1	B	673	ALA	3.8
1	A	705	GLU	3.5
1	B	614	GLU	3.5
1	B	723	GLU	3.4
1	B	573	GLU	3.3
1	B	670	ALA	3.2
1	A	671	VAL	3.2
1	B	615	GLN	3.2
1	B	707	THR	3.1
1	B	671	VAL	3.1
1	B	662	THR	3.1
1	B	605	ILE	3.0
1	A	642	GLU	3.0
1	A	516	PHE	2.8
1	A	518	LEU	2.8
1	B	669	TRP	2.8
1	A	670	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	674	MET	2.8
1	B	622	SER	2.6
1	B	580	TRP	2.6
1	B	642	GLU	2.6
1	A	673	ALA	2.5
1	A	636	LEU	2.5
1	B	668	LEU	2.4
1	A	674	MET	2.4
1	A	547	VAL	2.3
1	B	643	CYS	2.3
1	B	661	PHE	2.3
1	B	578	LEU	2.3
1	B	613	GLY	2.3
1	B	483	VAL	2.3
1	B	561	VAL	2.2
1	B	485	ALA	2.2
1	B	691	GLU	2.1
1	A	644	GLU	2.1
1	B	539	GLN	2.1
1	B	603	ALA	2.1
1	B	718	ILE	2.1
1	A	637	LYS	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 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.