



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:55 PM GMT

PDB ID : 3VH3  
Title : Crystal structure of Atg7CTD-Atg8 complex  
Authors : Noda, N.N.; Satoo, K.; Inagaki, F.  
Deposited on : 2011-08-23  
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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

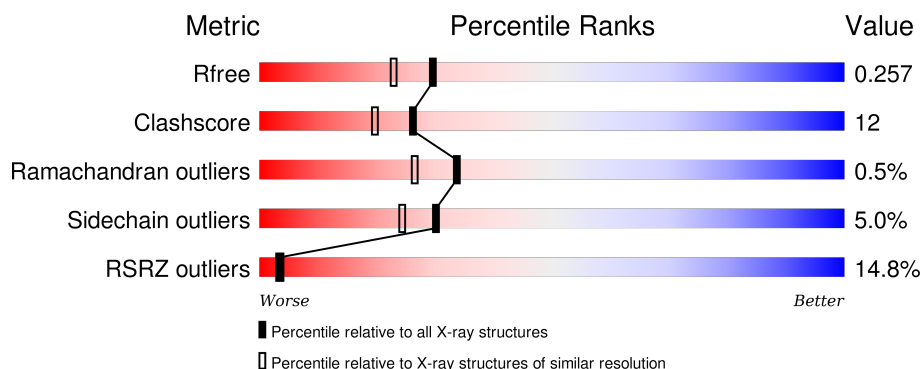
# 1 Overall quality at a glance

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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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	340	<div> <div>11%</div> <div>75%</div> <div>17%</div> <div>• 6%</div> </div>
2	B	119	<div> <div>22%</div> <div>60%</div> <div>34%</div> <div>• 5%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3459 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 Ubiquitin-like modifier-activating enzyme ATG7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2438	1560	415	448	15			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	291	GLY	-	EXPRESSION TAG	UNP P38862
A	292	PRO	-	EXPRESSION TAG	UNP P38862
A	293	HIS	-	EXPRESSION TAG	UNP P38862
A	294	MET	-	EXPRESSION TAG	UNP P38862

- Molecule 2 is a protein called Autophagy-related protein 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	113	Total	C	N	O	S	0	0	0
			852	557	137	155	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	EXPRESSION TAG	UNP P38182
B	-1	PRO	-	EXPRESSION TAG	UNP P38182
B	0	HIS	-	EXPRESSION TAG	UNP P38182
B	26	PRO	LYS	ENGINEERED MUTATION	UNP P38182

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

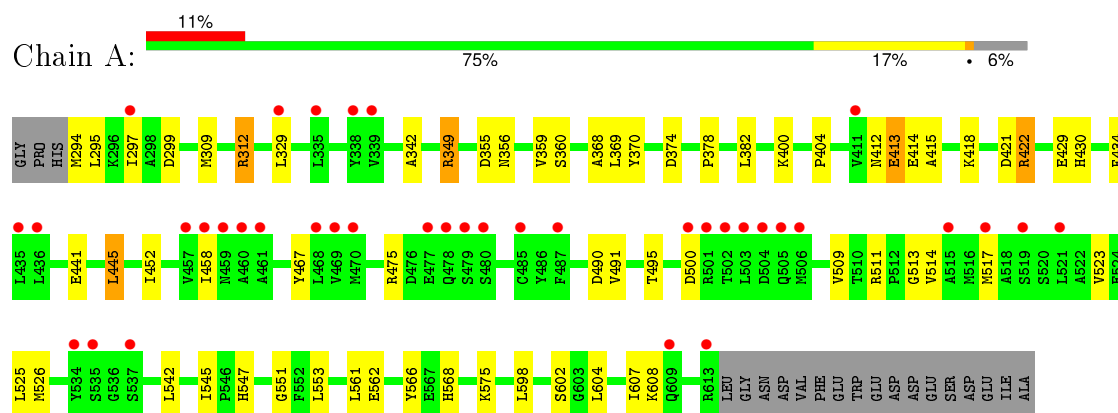
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	140	Total 140	O 140	0	0
4	B	28	Total 28	O 28	0	0

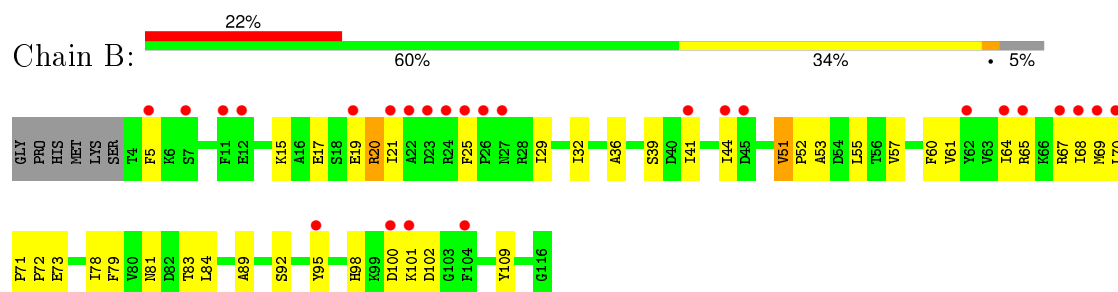
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Ubiquitin-like modifier-activating enzyme ATG7



- Molecule 2: Autophagy-related protein 8



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.81Å 69.81Å 220.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.28 – 2.00 33.28 – 2.00	Depositor EDS
% Data completeness (in resolution range)	91.6 (33.28-2.00) 91.6 (33.28-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.30 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.232 , 0.257 0.232 , 0.257	Depositor DCC
$R_{free}$ test set	3438 reflections (9.90%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.8	Xtriage
Anisotropy	0.599	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 55.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 37843 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3459	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.34	0/2486	0.61	0/3379
2	B	0.32	0/874	0.53	0/1191
All	All	0.34	0/3360	0.60	0/4570

There are no bond length outliers.

There are no bond angle outliers.

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	2438	0	2415	40	0
2	B	852	0	797	41	0
3	A	1	0	0	0	0
4	A	140	0	0	3	0
4	B	28	0	0	0	0
All	All	3459	0	3212	80	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 12.

All (80) 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:404:PRO:HG2	1:A:445:LEU:HD12	1.57	0.86
2:B:53:ALA:HB1	2:B:92:SER:HB3	1.59	0.82
2:B:70:LEU:HD12	2:B:71:PRO:HD2	1.60	0.81
1:A:349:ARG:HH11	1:A:349:ARG:HG2	1.46	0.79
2:B:60:PHE:CE2	2:B:64:ILE:HD11	2.21	0.76
2:B:71:PRO:HB2	2:B:73:GLU:OE1	1.86	0.75
1:A:561:LEU:HD13	2:B:84:LEU:HD13	1.73	0.70
2:B:52:PRO:HG2	2:B:55:LEU:HB2	1.75	0.69
1:A:374:ASP:HB3	1:A:382:LEU:HD11	1.79	0.64
1:A:349:ARG:HG2	1:A:349:ARG:NH1	2.11	0.64
2:B:29:ILE:HD11	2:B:53:ALA:HA	1.82	0.61
1:A:441:GLU:CG	1:A:495:THR:O	2.49	0.61
2:B:79:PHE:CE1	2:B:84:LEU:HD12	2.37	0.59
1:A:604:LEU:O	1:A:608:LYS:HG3	2.03	0.59
2:B:5:PHE:CD1	2:B:32:ILE:HD13	2.38	0.58
1:A:441:GLU:HG3	1:A:495:THR:O	2.03	0.58
1:A:370:TYR:HD2	1:A:382:LEU:HD22	1.70	0.57
2:B:17:GLU:O	2:B:21:ILE:HG13	2.05	0.56
2:B:81:ASN:ND2	2:B:98:HIS:HB3	2.21	0.56
1:A:509:VAL:HA	4:A:815:HOH:O	2.04	0.56
1:A:356:ASN:HA	1:A:400:LYS:HG3	1.87	0.56
2:B:20:ARG:HG2	2:B:20:ARG:HH11	1.72	0.55
2:B:64:ILE:HG22	2:B:68:ILE:HD12	1.89	0.54
2:B:57:VAL:HG23	2:B:89:ALA:O	2.08	0.54
1:A:452:ILE:HA	1:A:475:ARG:HH21	1.73	0.54
1:A:412:ASN:HB3	1:A:415:ALA:HB3	1.91	0.52
2:B:79:PHE:HE1	2:B:84:LEU:HD12	1.75	0.52
1:A:517:MET:HE3	1:A:553:LEU:HD11	1.92	0.51
2:B:36:ALA:HB3	2:B:39:SER:HB2	1.91	0.51
1:A:413:GLU:HG3	1:A:414:GLU:N	2.24	0.51
2:B:68:ILE:O	2:B:68:ILE:HG22	2.11	0.51
1:A:404:PRO:HG2	1:A:445:LEU:CD1	2.37	0.51
2:B:61:VAL:O	2:B:65:ARG:HG2	2.11	0.50
1:A:370:TYR:CD2	1:A:382:LEU:HD22	2.47	0.49
2:B:65:ARG:NH2	2:B:72:PRO:HA	2.28	0.49
2:B:29:ILE:N	2:B:51:VAL:O	2.46	0.48
2:B:57:VAL:O	2:B:61:VAL:HG23	2.13	0.48
1:A:458:ILE:CD1	1:A:526:MET:HB2	2.43	0.48
2:B:20:ARG:HG2	2:B:20:ARG:NH1	2.28	0.48
2:B:41:ILE:HD11	2:B:109:TYR:CD1	2.49	0.48
2:B:70:LEU:CD1	2:B:71:PRO:HD2	2.38	0.47
2:B:65:ARG:HH21	2:B:72:PRO:HA	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:41:ILE:HD11	2:B:109:TYR:HD1	1.79	0.47
2:B:64:ILE:O	2:B:68:ILE:HG13	2.15	0.47
1:A:490:ASP:OD1	1:A:491:VAL:N	2.39	0.47
1:A:418:LYS:O	1:A:422:ARG:HB2	2.15	0.47
2:B:29:ILE:CD1	2:B:53:ALA:HA	2.45	0.46
2:B:60:PHE:O	2:B:64:ILE:HG13	2.15	0.46
1:A:378:PRO:HB2	4:A:820:HOH:O	2.14	0.46
2:B:17:GLU:HG3	2:B:20:ARG:NH1	2.30	0.46
1:A:429:GLU:HG3	1:A:430:HIS:ND1	2.31	0.46
2:B:73:GLU:CD	2:B:73:GLU:H	2.20	0.45
1:A:607:ILE:HD12	1:A:608:LYS:N	2.32	0.45
1:A:547:HIS:HD2	1:A:562:GLU:HA	1.82	0.45
2:B:72:PRO:HG2	2:B:73:GLU:OE2	2.17	0.44
1:A:517:MET:CE	1:A:553:LEU:HD11	2.48	0.44
1:A:342:ALA:HB3	1:A:523:VAL:HG21	1.99	0.44
2:B:57:VAL:CG1	2:B:78:ILE:HD11	2.48	0.43
1:A:467:TYR:CZ	1:A:551:GLY:HA3	2.53	0.43
1:A:359:VAL:HG21	1:A:382:LEU:HD13	1.99	0.43
1:A:355:ASP:O	1:A:400:LYS:HA	2.17	0.43
1:A:312:ARG:HG2	1:A:312:ARG:HH11	1.82	0.43
1:A:299:ASP:OD1	1:A:349:ARG:NH2	2.52	0.43
1:A:607:ILE:HD12	1:A:607:ILE:C	2.38	0.43
2:B:17:GLU:HA	2:B:20:ARG:HG2	2.01	0.43
2:B:44:ILE:HD12	2:B:67:ARG:O	2.19	0.43
2:B:29:ILE:HG23	2:B:95:TYR:CD2	2.54	0.42
1:A:545:ILE:O	1:A:545:ILE:HG13	2.19	0.42
2:B:15:LYS:O	2:B:19:GLU:HG3	2.19	0.42
1:A:511:ARG:O	1:A:514:VAL:HG22	2.19	0.42
1:A:566:TYR:CE2	1:A:568:HIS:HB2	2.55	0.41
1:A:513:GLY:O	1:A:517:MET:HG3	2.21	0.41
1:A:294:MET:O	1:A:297:ILE:HG22	2.20	0.41
2:B:100:ASP:HB3	2:B:102:ASP:OD1	2.20	0.41
2:B:64:ILE:CG2	2:B:68:ILE:HD12	2.51	0.41
1:A:329:LEU:HD23	1:A:434:PHE:HB2	2.03	0.41
1:A:575:LYS:HE2	1:A:602:SER:O	2.20	0.41
1:A:514:VAL:HG12	1:A:553:LEU:HD13	2.03	0.40
2:B:100:ASP:OD1	2:B:101:LYS:N	2.55	0.40
4:A:905:HOH:O	2:B:83:THR:HG23	2.21	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	318/340 (94%)	307 (96%)	9 (3%)	2 (1%)	30	22
2	B	111/119 (93%)	105 (95%)	6 (5%)	0	100	100
All	All	429/459 (94%)	412 (96%)	15 (4%)	2 (0%)	34	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	368	ALA
1	A	413	GLU

### 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	259/298 (87%)	246 (95%)	13 (5%)	30	24
2	B	82/106 (77%)	78 (95%)	4 (5%)	31	25
All	All	341/404 (84%)	324 (95%)	17 (5%)	30	24

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	295	LEU
1	A	309	MET
1	A	312	ARG
1	A	349	ARG

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Mol	Chain	Res	Type
1	A	360	SER
1	A	369	LEU
1	A	421	ASP
1	A	422	ARG
1	A	445	LEU
1	A	500	ASP
1	A	525	LEU
1	A	542	LEU
1	A	598	LEU
2	B	20	ARG
2	B	25	PHE
2	B	51	VAL
2	B	69	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	417	HIS
1	A	547	HIS
2	B	81	ASN

### 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](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	320/340 (94%)	0.55	38 (11%) 6 6	28, 43, 74, 85	0
2	B	113/119 (94%)	1.25	26 (23%) 1 1	33, 71, 91, 96	0
All	All	433/459 (94%)	0.73	64 (14%) 3 3	28, 47, 84, 96	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	25	PHE	6.8
1	A	478	GLN	5.0
2	B	23	ASP	4.8
2	B	27	ASN	4.8
2	B	24	ARG	4.6
1	A	506	MET	4.2
2	B	44	ILE	4.1
1	A	460	ALA	4.0
1	A	339	VAL	4.0
2	B	5	PHE	3.9
1	A	534	TYR	3.8
2	B	104	PHE	3.8
2	B	62	TYR	3.8
2	B	22	ALA	3.7
1	A	477	GLU	3.7
1	A	436	LEU	3.7
2	B	21	ILE	3.5
1	A	505	GLN	3.4
1	A	609	GLN	3.3
2	B	19	GLU	3.3
2	B	45	ASP	3.2
1	A	458	ILE	3.2
1	A	338	TYR	3.1
2	B	69	MET	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	435	LEU	2.9
1	A	479	SER	2.8
1	A	501	ARG	2.8
1	A	461	ALA	2.8
1	A	504	ASP	2.8
2	B	100	ASP	2.8
2	B	11	PHE	2.8
2	B	41	ILE	2.7
1	A	515	ALA	2.7
1	A	487	PHE	2.7
2	B	68	ILE	2.7
1	A	519	SER	2.7
2	B	26	PRO	2.6
2	B	64	ILE	2.6
1	A	335	LEU	2.6
1	A	537	SER	2.6
1	A	503	LEU	2.6
2	B	67	ARG	2.6
1	A	500	ASP	2.6
1	A	468	LEU	2.5
1	A	459	ASN	2.5
2	B	101	LYS	2.4
1	A	457	VAL	2.4
1	A	411	VAL	2.4
2	B	70	LEU	2.4
1	A	535	SER	2.3
1	A	469	VAL	2.3
1	A	297	ILE	2.3
2	B	95	TYR	2.3
2	B	12	GLU	2.2
1	A	470	MET	2.2
2	B	7	SER	2.2
1	A	502	THR	2.2
1	A	480	SER	2.1
1	A	613	ARG	2.1
1	A	521	LEU	2.1
2	B	65	ARG	2.1
1	A	329	LEU	2.0
1	A	485	CYS	2.0
1	A	517	MET	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	A	701	1/1	0.98	0.11	-1.14	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.