



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 09:50 AM GMT

PDB ID : 3JW0  
Title : E2 Ubiquitin-HECT  
Authors : Kamadurai, H.B.; Schulman, B.A.  
Deposited on : 2009-09-17  
Resolution : 3.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

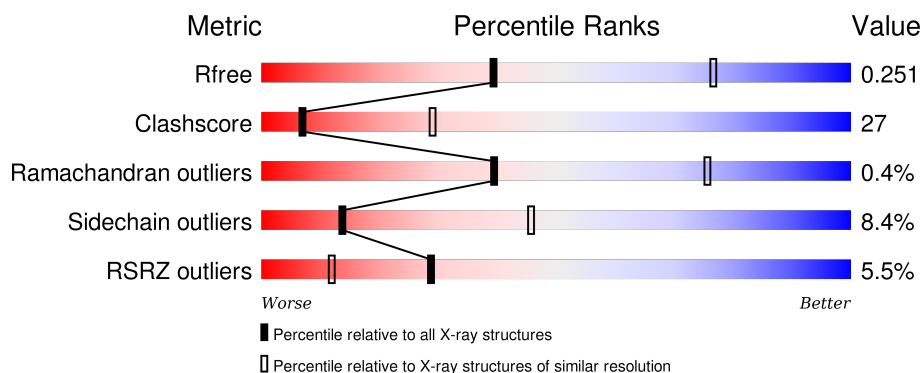
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	146	<div> <div>2%</div> <div> <div></div> <div>54%</div> <div>44%</div> <div>•</div> </div> </div>
1	B	146	<div> <div>%</div> <div> <div></div> <div>53%</div> <div>40%</div> <div>7%</div> </div> </div>
2	C	385	<div> <div>%</div> <div> <div></div> <div>49%</div> <div>45%</div> <div>• •</div> </div> </div>
2	D	385	<div> <div></div> <div> <div></div> <div>55%</div> <div>37%</div> <div>5%</div> <div>•</div> </div> </div>
3	X	81	<div> <div>38%</div> <div> <div></div> <div>40%</div> <div>53%</div> <div>• 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	Y	81	<div><div></div><div>33%</div><div>46%</div><div>46%</div><div>6%</div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9826 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-conjugating enzyme E2 D2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	0	0
			1170	749	202	213	6			
1	B	146	Total	C	N	O	S	0	0	0
			1170	749	202	213	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	SER	LEU	ENGINEERED	UNP P62837
A	85	SER	CYS	ENGINEERED	UNP P62837
A	98	LYS	THR	ENGINEERED	UNP P62837
B	3	SER	LEU	ENGINEERED	UNP P62837
B	85	SER	CYS	ENGINEERED	UNP P62837
B	98	LYS	THR	ENGINEERED	UNP P62837

- Molecule 2 is a protein called E3 ubiquitin-protein ligase NEDD4-like.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	376	Total	C	N	O	S	0	0	0
			3140	2032	519	572	17			
2	D	374	Total	C	N	O	S	0	0	0
			3125	2021	517	570	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	571	GLY	-	EXPRESSION TAG	UNP Q96PU5
C	572	SER	-	EXPRESSION TAG	UNP Q96PU5
C	573	PRO	-	EXPRESSION TAG	UNP Q96PU5
C	574	GLU	-	EXPRESSION TAG	UNP Q96PU5
C	575	PHE	-	EXPRESSION TAG	UNP Q96PU5
C	922	SER	CYS	ENGINEERED	UNP Q96PU5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	571	GLY	-	EXPRESSION TAG	UNP Q96PU5
D	572	SER	-	EXPRESSION TAG	UNP Q96PU5
D	573	PRO	-	EXPRESSION TAG	UNP Q96PU5
D	574	GLU	-	EXPRESSION TAG	UNP Q96PU5
D	575	PHE	-	EXPRESSION TAG	UNP Q96PU5
D	922	SER	CYS	ENGINEERED	UNP Q96PU5

- Molecule 3 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	X	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
3	Y	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	-4	GLY	-	EXPRESSION TAG	UNP P62988
X	-3	SER	-	EXPRESSION TAG	UNP P62988
X	-2	GLY	-	EXPRESSION TAG	UNP P62988
X	-1	GLY	-	EXPRESSION TAG	UNP P62988
X	0	SER	-	EXPRESSION TAG	UNP P62988
Y	-4	GLY	-	EXPRESSION TAG	UNP P62988
Y	-3	SER	-	EXPRESSION TAG	UNP P62988
Y	-2	GLY	-	EXPRESSION TAG	UNP P62988
Y	-1	GLY	-	EXPRESSION TAG	UNP P62988
Y	0	SER	-	EXPRESSION TAG	UNP P62988

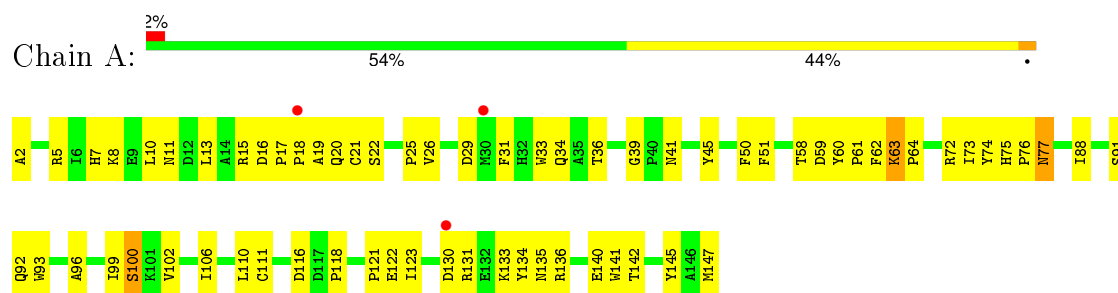
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O	0	0
			1	1		
4	C	6	Total	O	0	0
			6	6		
4	D	11	Total	O	0	0
			11	11		
4	Y	1	Total	O	0	0
			1	1		

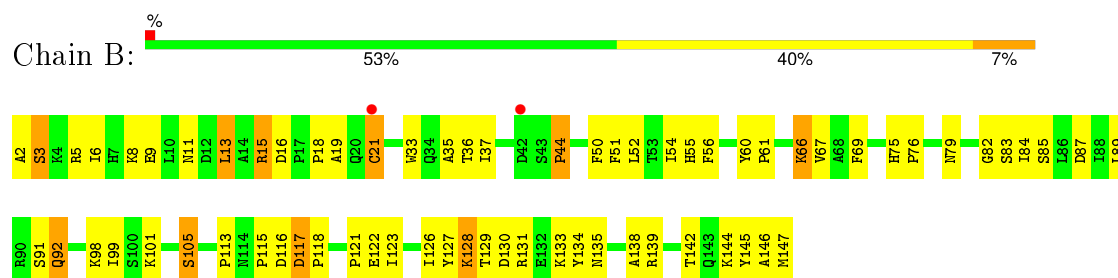
### 3 Residue-property plots [i](#)

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.

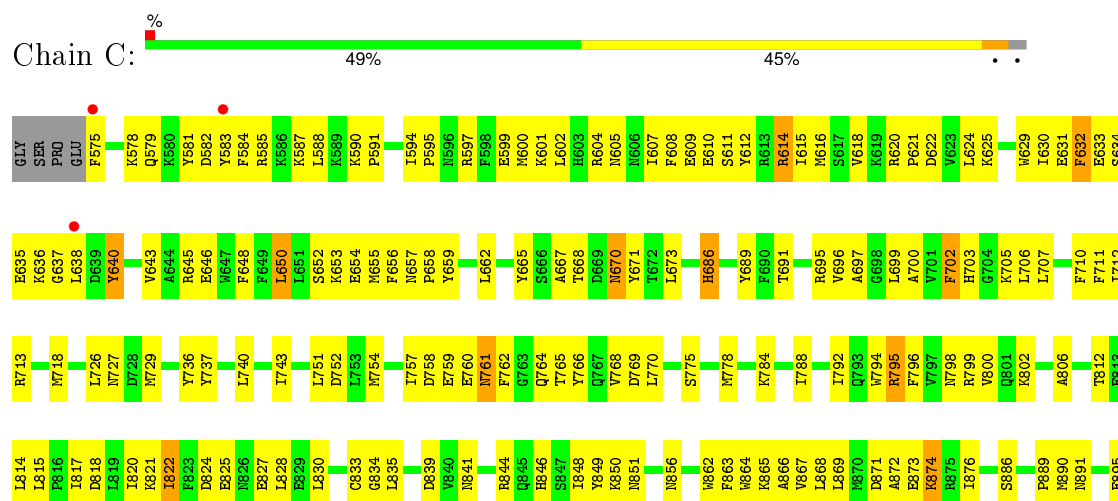
#### • Molecule 1: Ubiquitin-conjugating enzyme E2 D2



#### • Molecule 1: Ubiquitin-conjugating enzyme E2 D2

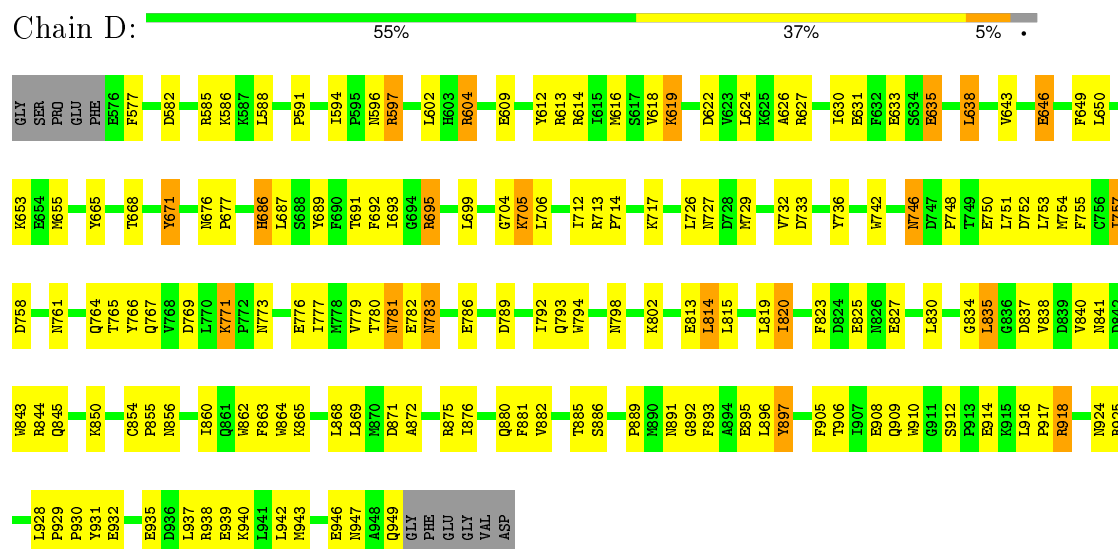


#### • Molecule 2: E3 ubiquitin-protein ligase NEDD4-like

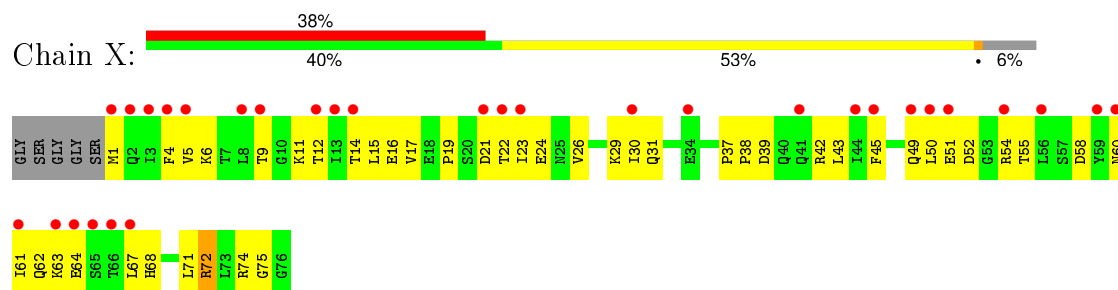




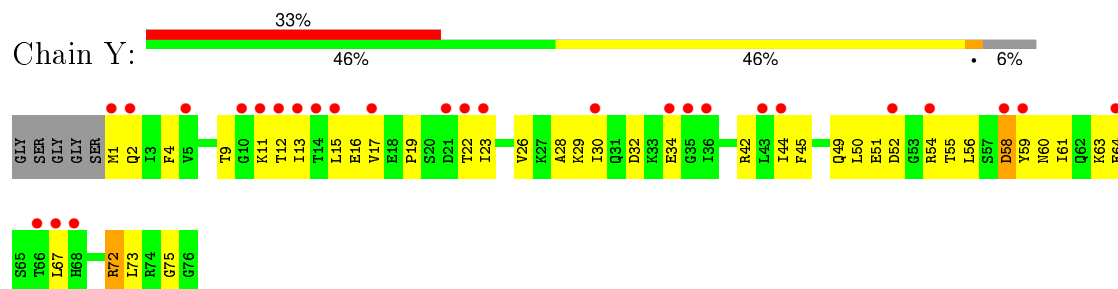
• Molecule 2: E3 ubiquitin-protein ligase NEDD4-like



• Molecule 3: Ubiquitin



• Molecule 3: Ubiquitin



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	174.17Å 200.57Å 109.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.10 49.72 – 3.12	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.10) 97.5 (49.72-3.12)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.46 (at 3.12Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.252 , 0.287 0.253 , 0.251	Depositor DCC
$R_{free}$ test set	1700 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	80.9	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 61.4	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	1 of 33799 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	9826	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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 [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	0.41	0/1206	0.70	0/1642
1	B	0.41	0/1206	0.66	0/1642
2	C	0.47	0/3224	0.68	0/4355
2	D	0.49	0/3208	0.71	0/4334
3	X	0.44	0/607	0.68	0/816
3	Y	0.38	0/607	0.61	0/816
All	All	0.46	0/10058	0.69	0/13605

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 [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	1170	0	1152	73	0
1	B	1170	0	1152	62	0
2	C	3140	0	3057	171	0
2	D	3125	0	3045	163	0
3	X	601	0	629	45	0
3	Y	601	0	629	31	0
4	B	1	0	0	0	0
4	C	6	0	0	1	0
4	D	11	0	0	2	0
4	Y	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9826	0	9664	518	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 27.

The worst 5 of 518 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:51:PHE:CE2	2:D:868:LEU:HD11	1.78	1.17
3:X:4:PHE:CD1	3:X:14:THR:HG22	1.95	1.01
2:C:906:THR:HB	2:C:925:ARG:HG3	1.47	0.96
2:D:733:ASP:OD2	2:D:736:TYR:HB2	1.69	0.93
1:B:116:ASP:O	1:B:118:PRO:HD3	1.69	0.92

There are no symmetry-related clashes.

## 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	144/146 (99%)	130 (90%)	14 (10%)	0	100	100
1	B	144/146 (99%)	132 (92%)	11 (8%)	1 (1%)	26	65
2	C	374/385 (97%)	310 (83%)	61 (16%)	3 (1%)	24	63
2	D	372/385 (97%)	315 (85%)	56 (15%)	1 (0%)	46	80
3	X	74/81 (91%)	66 (89%)	8 (11%)	0	100	100
3	Y	74/81 (91%)	65 (88%)	9 (12%)	0	100	100
All	All	1182/1224 (97%)	1018 (86%)	159 (14%)	5 (0%)	39	75

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	686	HIS
2	D	881	PHE
2	C	595	PRO
2	C	800	VAL
1	B	44	PRO

### 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	130/130 (100%)	123 (95%)	7 (5%)	27 64
1	B	130/130 (100%)	117 (90%)	13 (10%)	9 34
2	C	341/348 (98%)	315 (92%)	26 (8%)	16 51
2	D	340/348 (98%)	305 (90%)	35 (10%)	9 32
3	X	68/70 (97%)	64 (94%)	4 (6%)	24 60
3	Y	68/70 (97%)	63 (93%)	5 (7%)	17 51
All	All	1077/1096 (98%)	987 (92%)	90 (8%)	14 46

5 of 90 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	841	ASN
2	D	619	LYS
3	X	58	ASP
2	C	868	LEU
2	C	943	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	798	ASN
2	D	596	ASN
3	X	62	GLN
2	C	861	GLN

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Mol	Chain	Res	Type
2	D	676	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](#)

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	146/146 (100%)	0.27	3 (2%) 67 44	74, 102, 110, 110	0
1	B	146/146 (100%)	0.30	2 (1%) 78 60	73, 98, 110, 110	0
2	C	376/385 (97%)	0.23	3 (0%) 87 75	64, 96, 110, 110	0
2	D	374/385 (97%)	0.14	0 100 100	63, 90, 110, 110	0
3	X	76/81 (93%)	1.90	31 (40%) 0 0	107, 110, 110, 110	0
3	Y	76/81 (93%)	1.64	27 (35%) 0 0	105, 110, 110, 110	0
All	All	1194/1224 (97%)	0.41	66 (5%) 29 12	63, 98, 110, 110	0

The worst 5 of 66 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	X	65	SER	6.3
3	X	66	THR	6.0
3	X	23	ILE	5.3
3	X	3	ILE	5.0
3	X	2	GLN	4.9

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.