



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:52 pm GMT

PDB ID : 1ZMW  
Title : Catalytic and ubiquitin-associated domains of MARK2/PAR-1: T208A/S212A inactive double mutant  
Authors : Panneerselvam, S.; Marx, A.; Mandelkow, E.-M.; Mandelkow, E.  
Deposited on : 2005-05-11  
Resolution : 2.80 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  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

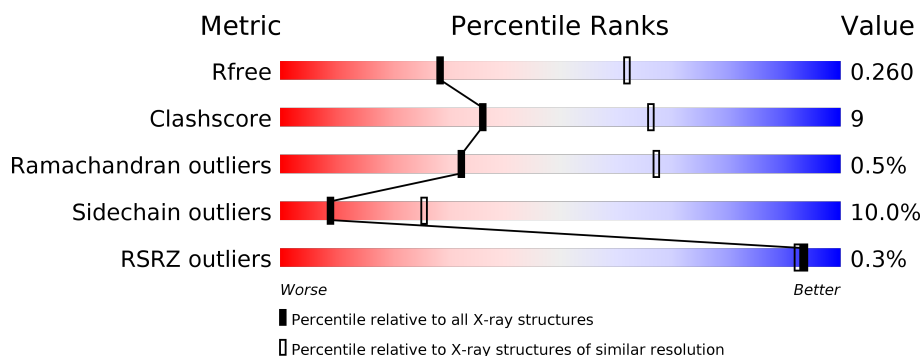
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	327	
1	B	327	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4885 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 MAP/Microtubule affinity regulating kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2437	1570	415	439	13			
1	B	302	Total	C	N	O	S	0	0	0
			2448	1574	421	440	13			

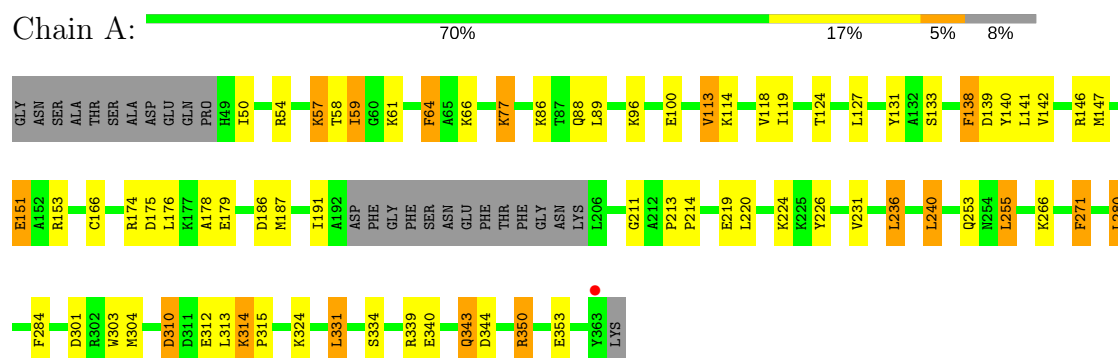
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	38	GLY	-	CLONING ARTIFACT	UNP O08679
A	208	ALA	THR	ENGINEERED	UNP O08679
A	212	ALA	SER	ENGINEERED	UNP O08679
B	38	GLY	-	CLONING ARTIFACT	UNP O08679
B	208	ALA	THR	ENGINEERED	UNP O08679
B	212	ALA	SER	ENGINEERED	UNP O08679

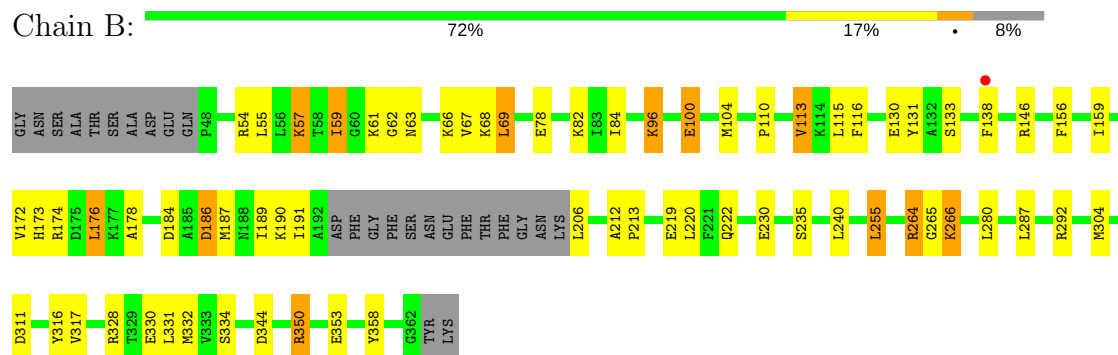
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: MAP/Microtubule affinity regulating kinase 2



- Molecule 1: MAP/Microtubule affinity regulating kinase 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.30Å 119.30Å 105.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	73.88 – 2.80 73.88 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (73.88-2.80) 100.0 (73.88-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.02 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, $R_{free}$	0.203 , 0.262 0.201 , 0.260	Depositor DCC
$R_{free}$ test set	1092 reflections (5.47%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.0	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 27.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.040 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4885	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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	1.02	3/2485 (0.1%)	0.94	3/3348 (0.1%)
1	B	0.98	2/2496 (0.1%)	0.94	3/3360 (0.1%)
All	All	1.00	5/4981 (0.1%)	0.94	6/6708 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	100	GLU	CB-CG	8.30	1.68	1.52
1	B	100	GLU	CG-CD	7.92	1.63	1.51
1	A	100	GLU	CG-CD	7.90	1.63	1.51
1	B	100	GLU	CB-CG	7.01	1.65	1.52
1	A	166	CYS	CB-SG	-5.90	1.72	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	ARG	NE-CZ-NH2	-7.51	116.54	120.30
1	B	328	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	B	62	GLY	N-CA-C	-5.41	99.57	113.10
1	A	59	ILE	CB-CA-C	-5.33	100.93	111.60
1	A	77	LYS	CD-CE-NZ	5.17	123.59	111.70
1	B	311	ASP	CB-CG-OD1	-5.03	113.78	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	310	ASP	Peptide

## 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	2437	0	2486	43	0
1	B	2448	0	2511	48	0
All	All	4885	0	4997	90	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 9.

All (90) 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:57:LYS:HE2	1:A:58:THR:H	1.07	1.11
1:A:59:ILE:HG22	1:A:59:ILE:O	1.62	0.98
1:A:280:LEU:HD11	1:A:304:MET:CE	1.99	0.93
1:A:57:LYS:HE2	1:A:58:THR:N	1.88	0.89
1:B:264:ARG:NH2	1:B:266:LYS:HE3	1.88	0.87
1:A:138:PHE:HB2	1:A:178:ALA:HB1	1.62	0.81
1:A:59:ILE:CG2	1:A:59:ILE:O	2.29	0.80
1:A:57:LYS:CE	1:A:58:THR:H	1.94	0.79
1:B:332:MET:CE	1:B:358:TYR:HD1	1.96	0.78
1:B:69:LEU:HD21	1:B:78:GLU:OE1	1.84	0.76
1:B:280:LEU:HD21	1:B:304:MET:CE	2.17	0.75
1:A:313:LEU:O	1:A:314:LYS:HE2	1.88	0.74
1:B:280:LEU:HD21	1:B:304:MET:HE2	1.70	0.72
1:A:314:LYS:HD3	1:A:315:PRO:HD2	1.72	0.71
1:B:96:LYS:O	1:B:100:GLU:HG3	1.94	0.68
1:A:280:LEU:HD11	1:A:304:MET:HE3	1.74	0.66
1:B:332:MET:HE1	1:B:358:TYR:HD1	1.58	0.65
1:A:113:VAL:HG13	1:A:191:ILE:O	1.98	0.63
1:A:50:ILE:HG22	1:A:119:ILE:HD12	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LYS:HE2	1:B:68:LYS:HE3	1.82	0.61
1:A:331:LEU:O	1:A:334:SER:HB2	1.99	0.61
1:A:186:ASP:O	1:A:187:MET:HB2	2.01	0.61
1:B:332:MET:HE2	1:B:358:TYR:HD1	1.66	0.60
1:B:82:LYS:HD2	1:B:84:ILE:HD11	1.83	0.59
1:B:55:LEU:HD13	1:B:68:LYS:HD3	1.85	0.57
1:B:264:ARG:HH21	1:B:266:LYS:HE3	1.70	0.57
1:A:131:TYR:CE2	1:A:133:SER:HB3	2.39	0.57
1:B:332:MET:HE2	1:B:358:TYR:CD1	2.39	0.56
1:B:331:LEU:O	1:B:334:SER:HB2	2.05	0.56
1:B:69:LEU:CD2	1:B:78:GLU:OE1	2.55	0.55
1:A:280:LEU:HD12	1:A:303:TRP:CE3	2.41	0.54
1:B:156:PHE:CE2	1:B:304:MET:HE3	2.42	0.54
1:B:67:VAL:HG22	1:B:82:LYS:HG2	1.89	0.54
1:B:172:VAL:HG23	1:B:174:ARG:HG3	1.89	0.53
1:A:343:GLN:NE2	1:A:344:ASP:OD1	2.41	0.53
1:B:173:HIS:O	1:B:174:ARG:HB2	2.09	0.53
1:A:138:PHE:CD1	1:A:138:PHE:C	2.83	0.52
1:B:186:ASP:O	1:B:187:MET:HB2	2.10	0.52
1:B:113:VAL:HG12	1:B:190:LYS:HB3	1.91	0.52
1:A:344:ASP:HB3	1:A:350:ARG:NH2	2.26	0.51
1:B:138:PHE:HB2	1:B:178:ALA:HB1	1.93	0.50
1:A:138:PHE:HD1	1:A:138:PHE:C	2.15	0.50
1:A:175:ASP:N	1:A:175:ASP:OD1	2.45	0.49
1:B:156:PHE:HD2	1:B:304:MET:HE1	1.78	0.49
1:A:240:LEU:HD11	1:A:280:LEU:HD13	1.93	0.49
1:A:64:PHE:CD2	1:A:88:GLN:HB3	2.48	0.49
1:B:159:ILE:HG12	1:B:189:ILE:HG21	1.94	0.49
1:A:220:LEU:HB2	1:A:226:TYR:CE1	2.48	0.48
1:A:280:LEU:HD22	1:A:284:PHE:CE2	2.48	0.48
1:B:104:MET:HG3	1:B:115:LEU:HB2	1.94	0.48
1:B:156:PHE:CD2	1:B:304:MET:CE	2.96	0.48
1:B:332:MET:CE	1:B:358:TYR:CD1	2.86	0.48
1:A:174:ARG:HD3	1:A:226:TYR:O	2.14	0.48
1:A:280:LEU:HG	1:A:301:ASP:OD2	2.13	0.48
1:A:271:PHE:C	1:A:271:PHE:CD1	2.87	0.47
1:B:230:GLU:N	1:B:230:GLU:OE1	2.43	0.47
1:B:344:ASP:HB3	1:B:350:ARG:NH2	2.30	0.46
1:B:219:GLU:OE2	1:B:292:ARG:NH2	2.49	0.46
1:A:139:ASP:O	1:A:140:TYR:C	2.54	0.45
1:B:212:ALA:HB3	1:B:213:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:LEU:HD21	1:B:304:MET:HE1	1.97	0.45
1:A:303:TRP:HE3	1:A:304:MET:HE2	1.80	0.45
1:B:156:PHE:HD2	1:B:304:MET:CE	2.29	0.45
1:B:113:VAL:HG13	1:B:191:ILE:O	2.17	0.45
1:A:140:TYR:C	1:A:140:TYR:CD1	2.91	0.44
1:A:213:PRO:N	1:A:214:PRO:CD	2.81	0.44
1:B:219:GLU:HA	1:B:222:GLN:HE21	1.83	0.44
1:A:211:GLY:O	1:A:214:PRO:HD2	2.18	0.44
1:A:224:LYS:HA	1:A:224:LYS:HD3	1.75	0.44
1:A:64:PHE:HD2	1:A:88:GLN:HB3	1.82	0.44
1:B:184:ASP:OD1	1:B:184:ASP:C	2.57	0.43
1:B:57:LYS:HD2	1:B:59:ILE:HG22	2.00	0.43
1:A:86:LYS:HA	1:A:89:LEU:HD12	2.01	0.43
1:A:236:LEU:HA	1:A:236:LEU:HD12	1.93	0.42
1:A:255:LEU:HD12	1:B:220:LEU:HD21	2.01	0.42
1:A:147:MET:HB3	1:A:151:GLU:HB3	2.00	0.42
1:A:219:GLU:HB2	1:A:224:LYS:HB2	2.00	0.42
1:A:255:LEU:N	1:A:255:LEU:HD23	2.34	0.42
1:B:265:GLY:HA3	1:B:287:LEU:HD21	2.02	0.41
1:B:110:PRO:HD3	1:B:316:TYR:OH	2.20	0.41
1:B:116:PHE:HE2	1:B:130:GLU:HA	1.85	0.41
1:B:176:LEU:HA	1:B:176:LEU:HD12	1.88	0.41
1:A:303:TRP:CE3	1:A:304:MET:HE2	2.55	0.41
1:A:139:ASP:O	1:A:142:VAL:HG12	2.20	0.41
1:B:156:PHE:CD2	1:B:304:MET:HE1	2.55	0.41
1:B:131:TYR:CE2	1:B:133:SER:HB3	2.56	0.41
1:B:156:PHE:CD2	1:B:304:MET:HE3	2.56	0.41
1:B:61:LYS:HE2	1:B:61:LYS:HB3	1.89	0.41
1:B:255:LEU:HD13	1:B:255:LEU:HA	1.84	0.40
1:B:280:LEU:CD2	1:B:304:MET:HE2	2.45	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	298/327 (91%)	278 (93%)	18 (6%)	2 (1%)	25	59
1	B	298/327 (91%)	283 (95%)	14 (5%)	1 (0%)	44	77
All	All	596/654 (91%)	561 (94%)	32 (5%)	3 (0%)	32	67

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146	ARG
1	A	179	GLU
1	B	146	ARG

### 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	264/287 (92%)	230 (87%)	34 (13%)	5	15
1	B	267/287 (93%)	248 (93%)	19 (7%)	17	44
All	All	531/574 (92%)	478 (90%)	53 (10%)	9	26

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	57	LYS
1	A	61	LYS
1	A	64	PHE
1	A	66	LYS
1	A	77	LYS
1	A	96	LYS
1	A	113	VAL
1	A	114	LYS
1	A	118	VAL

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Mol	Chain	Res	Type
1	A	124	THR
1	A	127	LEU
1	A	138	PHE
1	A	141	LEU
1	A	151	GLU
1	A	176	LEU
1	A	231	VAL
1	A	236	LEU
1	A	240	LEU
1	A	253	GLN
1	A	255	LEU
1	A	266	LYS
1	A	271	PHE
1	A	280	LEU
1	A	310	ASP
1	A	312	GLU
1	A	314	LYS
1	A	324	LYS
1	A	331	LEU
1	A	339	ARG
1	A	340	GLU
1	A	343	GLN
1	A	350	ARG
1	A	353	GLU
1	B	54	ARG
1	B	57	LYS
1	B	59	ILE
1	B	63	ASN
1	B	69	LEU
1	B	96	LYS
1	B	113	VAL
1	B	176	LEU
1	B	186	ASP
1	B	206	LEU
1	B	235	SER
1	B	240	LEU
1	B	255	LEU
1	B	264	ARG
1	B	266	LYS
1	B	317	VAL
1	B	330	GLU
1	B	350	ARG

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Mol	Chain	Res	Type
1	B	353	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	63	ASN
1	B	144	HIS
1	B	222	GLN
1	B	253	GLN
1	B	343	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 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	302/327 (92%)	-0.21	1 (0%) 93 92	31, 46, 62, 70	0
1	B	302/327 (92%)	-0.24	1 (0%) 93 92	31, 47, 61, 69	0
All	All	604/654 (92%)	-0.22	2 (0%) 93 92	31, 46, 62, 70	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	363	TYR	2.3
1	B	138	PHE	2.3

### 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.