



# Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 11:17 am BST

PDB ID : 1Q88  
Title : Crystal structure of the C-domain of the T.vaginalis Inr binding protein, IBP39 (monoclinic form)  
Authors : Schumacher, M.A.; Johnson, P.J.  
Deposited on : 2003-08-20  
Resolution : 2.42 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.

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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 : 2.11  
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) : 2.11

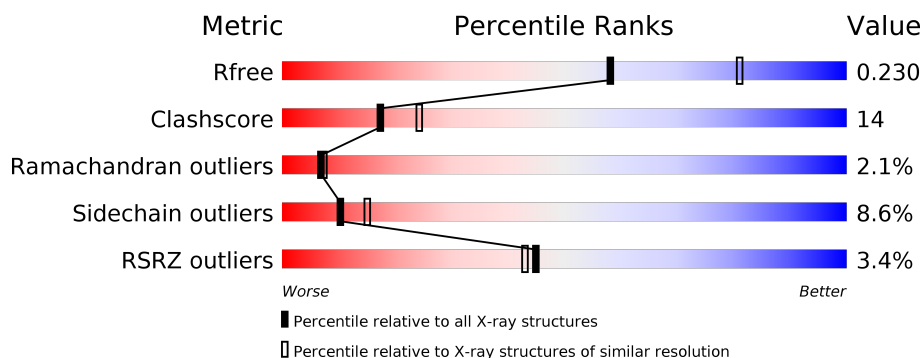
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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 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	221	<div> <div>4%</div> <div> <div></div> <div>59%</div> <div>21%</div> <div>5%</div> <div>14%</div> </div> </div>
1	B	221	<div> <div>2%</div> <div> <div></div> <div>62%</div> <div>24%</div> <div>•</div> <div>10%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3173 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 39 kDa initiator binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1540	988	251	292	9			
1	B	198	Total	C	N	O	S	0	0	0
			1603	1028	260	305	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	342	HIS	-	EXPRESSION TAG	UNP Q95VR4
A	343	HIS	-	EXPRESSION TAG	UNP Q95VR4
A	344	HIS	-	EXPRESSION TAG	UNP Q95VR4
A	345	HIS	-	EXPRESSION TAG	UNP Q95VR4
A	346	HIS	-	EXPRESSION TAG	UNP Q95VR4
A	347	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	342	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	343	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	344	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	345	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	346	HIS	-	EXPRESSION TAG	UNP Q95VR4
B	347	HIS	-	EXPRESSION TAG	UNP Q95VR4

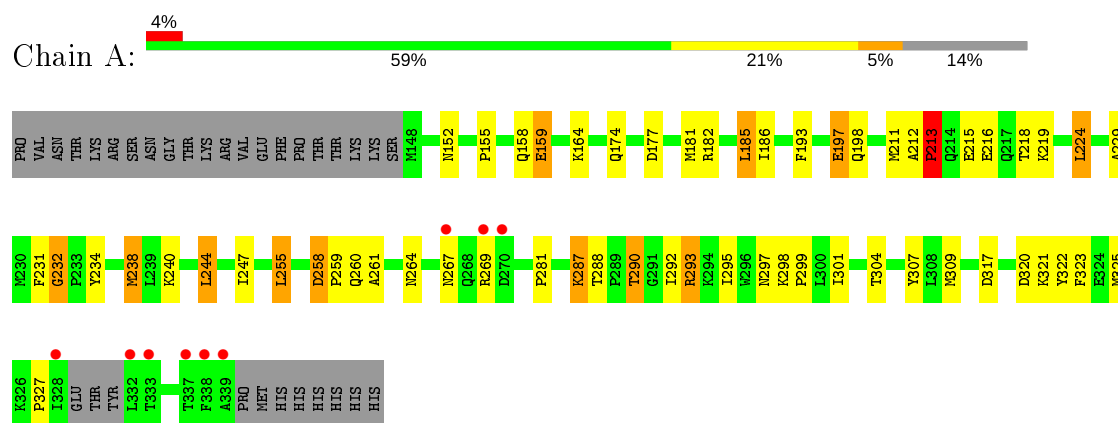
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	17	Total	O	0	0
			17	17		
2	B	13	Total	O	0	0
			13	13		

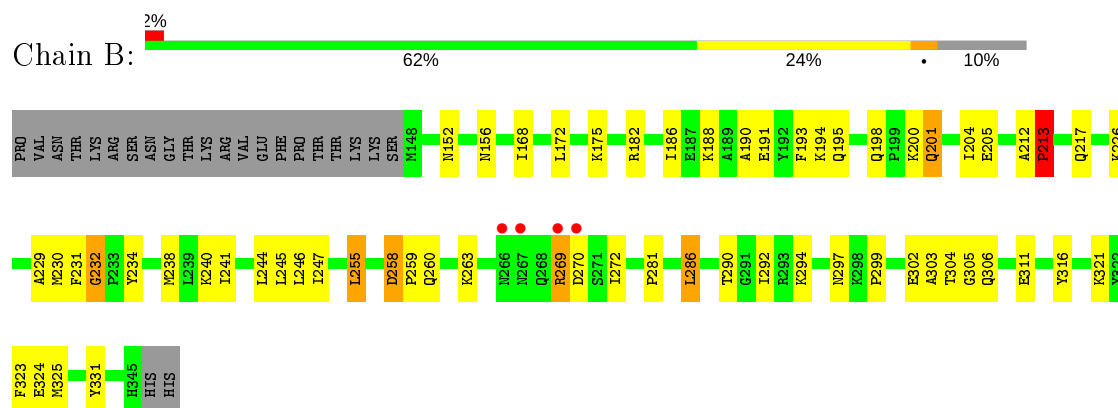
### 3 Residue-property plots [i](#)

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: 39 kDa initiator binding protein



- Molecule 1: 39 kDa initiator binding protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	99.60Å 76.40Å 80.30Å 90.00° 93.00° 90.00°	Depositor
Resolution (Å)	49.73 – 2.42 49.73 – 2.34	Depositor EDS
% Data completeness (in resolution range)	84.9 (49.73-2.42) 79.2 (49.73-2.34)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.19 (at 2.34Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.190 , 0.232 0.190 , 0.230	Depositor DCC
$R_{free}$ test set	1939 reflections (9.54%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.8	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 51.2	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	0.95	EDS
Total number of atoms	3173	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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	0.45	0/1577	0.83	5/2132 (0.2%)
1	B	0.48	0/1643	0.84	5/2225 (0.2%)
All	All	0.46	0/3220	0.83	10/4357 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	212	ALA	C-N-CD	-17.22	82.71	120.60
1	A	212	ALA	C-N-CD	-15.89	85.63	120.60
1	B	212	ALA	C-N-CA	13.74	179.71	122.00
1	A	212	ALA	C-N-CA	13.16	177.28	122.00
1	B	213	PRO	CA-N-CD	-10.36	97.00	111.50
1	A	213	PRO	CA-N-CD	-10.31	97.06	111.50
1	A	213	PRO	N-CD-CG	8.26	115.59	103.20
1	B	213	PRO	N-CD-CG	7.99	115.18	103.20
1	B	213	PRO	N-CA-CB	7.19	111.92	103.30
1	A	213	PRO	N-CA-CB	5.98	110.48	103.30

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	1540	0	1511	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1603	0	1558	44	0
2	A	17	0	0	1	0
2	B	13	0	0	2	0
All	All	3173	0	3069	86	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 14.

All (86) 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:B:198:GLN:HE22	1:B:238:MET:HG3	1.28	0.98
1:B:191:GLU:HG2	1:B:200:LYS:HD2	1.47	0.97
1:A:159:GLU:HG3	1:B:175:LYS:HA	1.52	0.90
1:B:269:ARG:HD3	1:B:269:ARG:H	1.53	0.73
1:A:232:GLY:H	1:A:297:ASN:HD21	1.38	0.72
1:A:297:ASN:O	1:A:299:PRO:HD3	1.91	0.69
1:A:182:ARG:HD3	2:A:361:HOH:O	1.95	0.67
1:B:195:GLN:H	1:B:198:GLN:HE21	1.43	0.65
1:B:260:GLN:HG3	1:B:263:LYS:HB2	1.80	0.63
1:A:317:ASP:H	1:A:321:LYS:HZ3	1.47	0.63
1:B:272:ILE:HD11	1:B:331:TYR:HB2	1.82	0.61
1:A:155:PRO:HA	1:A:158:GLN:HE21	1.66	0.61
1:A:159:GLU:CG	1:B:175:LYS:HA	2.30	0.60
1:A:264:ASN:HA	1:A:267:ASN:HD22	1.65	0.60
1:B:229:ALA:HB3	1:B:281:PRO:HB2	1.82	0.59
1:B:244:LEU:HD11	1:B:255:LEU:HG	1.84	0.59
1:A:186:ILE:HD13	1:A:211:MET:HE2	1.82	0.59
1:B:188:LYS:HE2	1:B:188:LYS:HA	1.84	0.59
1:B:297:ASN:O	1:B:299:PRO:HD3	2.04	0.57
1:B:232:GLY:HA2	1:B:299:PRO:HA	1.85	0.57
1:A:298:LYS:HB3	1:A:301:ILE:HD12	1.88	0.56
1:B:200:LYS:O	1:B:204:ILE:HG13	2.04	0.56
1:A:298:LYS:CB	1:A:301:ILE:HD12	2.36	0.56
1:B:232:GLY:HA3	1:B:240:LYS:NZ	2.22	0.55
1:A:264:ASN:HA	1:A:267:ASN:ND2	2.21	0.55
1:B:195:GLN:H	1:B:198:GLN:NE2	2.04	0.54
1:A:317:ASP:HB2	1:A:321:LYS:HD3	1.89	0.53
1:B:182:ARG:O	1:B:186:ILE:HG12	2.09	0.52
1:A:229:ALA:HB3	1:A:281:PRO:HB2	1.90	0.52
1:A:293:ARG:HH11	1:A:293:ARG:HB3	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:ALA:CB	1:B:281:PRO:HB2	2.39	0.52
1:A:232:GLY:HA2	1:A:299:PRO:HA	1.92	0.51
1:A:258:ASP:O	1:A:261:ALA:HB2	2.12	0.50
1:B:241:ILE:O	1:B:245:LEU:HB2	2.11	0.50
1:A:216:GLU:HG2	1:A:218:THR:O	2.12	0.49
1:B:302:GLU:C	1:B:304:THR:H	2.15	0.49
1:A:247:ILE:HD13	1:A:323:PHE:HD2	1.77	0.49
1:A:198:GLN:OE1	1:A:238:MET:HG2	2.13	0.48
1:B:324:GLU:HB3	2:B:359:HOH:O	2.14	0.48
1:A:244:LEU:HD13	1:A:255:LEU:HD11	1.96	0.48
1:B:294:LYS:HB2	1:B:311:GLU:HG2	1.95	0.48
1:A:181:MET:CE	1:A:218:THR:HG22	2.44	0.47
1:B:240:LYS:NZ	1:B:297:ASN:HD21	2.12	0.47
1:B:306:GLN:HA	1:B:316:TYR:O	2.14	0.47
1:B:201:GLN:O	1:B:205:GLU:HG3	2.15	0.47
1:B:193:PHE:CE1	1:B:234:TYR:HB2	2.50	0.47
1:A:325:MET:C	1:A:327:PRO:HD3	2.34	0.47
1:A:247:ILE:HD11	1:A:320:ASP:HA	1.97	0.47
1:B:294:LYS:HB2	1:B:311:GLU:CG	2.45	0.47
1:B:226:LYS:O	1:B:230:MET:HG3	2.14	0.46
1:A:182:ARG:O	1:A:186:ILE:HG12	2.15	0.46
1:B:232:GLY:HA3	1:B:240:LYS:HZ2	1.80	0.46
1:A:293:ARG:CB	1:A:293:ARG:HH11	2.29	0.46
1:B:231:PHE:O	1:B:232:GLY:O	2.33	0.46
1:A:240:LYS:HZ3	1:A:307:TYR:HB2	1.80	0.45
1:B:232:GLY:CA	1:B:240:LYS:HZ2	2.28	0.45
1:A:185:LEU:HD13	1:A:211:MET:HE1	1.98	0.45
1:A:298:LYS:CE	1:A:309:MET:HE2	2.46	0.45
1:A:240:LYS:NZ	1:A:307:TYR:HB2	2.33	0.44
1:B:321:LYS:O	1:B:325:MET:HG2	2.17	0.44
1:A:298:LYS:HD3	1:A:309:MET:HE2	1.99	0.44
1:A:197:GLU:CD	1:A:197:GLU:H	2.21	0.43
1:A:224:LEU:HD23	1:A:224:LEU:HA	1.88	0.43
1:A:231:PHE:O	1:A:232:GLY:O	2.37	0.43
1:A:193:PHE:CD1	1:A:234:TYR:HB2	2.54	0.43
1:A:287:LYS:HG2	1:A:292:ILE:CD1	2.48	0.43
1:A:317:ASP:H	1:A:321:LYS:NZ	2.15	0.43
1:B:168:ILE:O	1:B:172:LEU:HG	2.19	0.43
1:A:232:GLY:H	1:A:297:ASN:ND2	2.11	0.42
1:A:232:GLY:N	1:A:297:ASN:HD21	2.12	0.42
1:B:292:ILE:HG13	1:B:292:ILE:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:GLN:NE2	1:B:238:MET:HG3	2.12	0.42
1:B:258:ASP:HA	1:B:259:PRO:HD3	1.87	0.42
1:B:190:ALA:O	1:B:194:LYS:HG2	2.20	0.41
1:B:246:LEU:HD23	2:B:360:HOH:O	2.20	0.41
1:A:295:ILE:HD13	1:A:322:TYR:CE2	2.56	0.41
1:B:290:THR:HG22	1:B:290:THR:O	2.20	0.41
1:A:182:ARG:HD2	1:A:216:GLU:O	2.20	0.41
1:A:288:THR:OG1	1:A:290:THR:HG23	2.20	0.41
1:B:302:GLU:C	1:B:304:THR:N	2.74	0.41
1:B:247:ILE:HD13	1:B:323:PHE:HD2	1.86	0.41
1:B:240:LYS:HZ2	1:B:297:ASN:HD21	1.68	0.41
1:A:181:MET:HE1	1:A:218:THR:HG22	2.03	0.40
1:A:258:ASP:HA	1:A:259:PRO:HD3	1.93	0.40
1:B:193:PHE:CD1	1:B:234:TYR:HB2	2.57	0.40
1:B:286:LEU:HD23	1:B:286:LEU:HA	1.87	0.40

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	185/221 (84%)	168 (91%)	13 (7%)	4 (2%)	6	7
1	B	196/221 (89%)	179 (91%)	13 (7%)	4 (2%)	7	8
All	All	381/442 (86%)	347 (91%)	26 (7%)	8 (2%)	7	7

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	232	GLY
1	B	213	PRO
1	B	232	GLY

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Mol	Chain	Res	Type
1	A	269	ARG
1	B	303	ALA
1	A	177	ASP
1	B	305	GLY
1	A	213	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	167/198 (84%)	148 (89%)	19 (11%)	5	7
1	B	172/198 (87%)	162 (94%)	10 (6%)	20	31
All	All	339/396 (86%)	310 (91%)	29 (9%)	10	15

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	159	GLU
1	A	164	LYS
1	A	174	GLN
1	A	185	LEU
1	A	197	GLU
1	A	213	PRO
1	A	215	GLU
1	A	219	LYS
1	A	224	LEU
1	A	238	MET
1	A	244	LEU
1	A	255	LEU
1	A	258	ASP
1	A	260	GLN
1	A	287	LYS
1	A	290	THR
1	A	293	ARG

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Mol	Chain	Res	Type
1	A	304	THR
1	B	152	ASN
1	B	156	ASN
1	B	201	GLN
1	B	213	PRO
1	B	217	GLN
1	B	255	LEU
1	B	258	ASP
1	B	269	ARG
1	B	270	ASP
1	B	286	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	214	GLN
1	A	260	GLN
1	A	267	ASN
1	A	297	ASN
1	A	312	ASN
1	B	152	ASN
1	B	198	GLN
1	B	214	GLN
1	B	253	HIS
1	B	306	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

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 [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	189/221 (85%)	-0.01	9 (4%) 30 28	31, 45, 112, 123	0
1	B	198/221 (89%)	-0.12	4 (2%) 65 62	31, 56, 100, 135	0
All	All	387/442 (87%)	-0.06	13 (3%) 45 43	31, 50, 108, 135	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	339	ALA	7.3
1	A	338	PHE	6.3
1	A	337	THR	5.9
1	B	270	ASP	5.5
1	B	269	ARG	4.7
1	A	332	LEU	3.6
1	A	333	THR	3.4
1	B	267	ASN	3.2
1	A	270	ASP	3.2
1	B	266	ASN	3.1
1	A	269	ARG	2.8
1	A	328	ILE	2.1
1	A	267	ASN	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.