



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 12:11 pm GMT

PDB ID : 1Q1T  
Title : Mouse Importin alpha: non-phosphorylated SV40 CN peptide complex  
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Deposited on : 2003-07-22  
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)

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	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

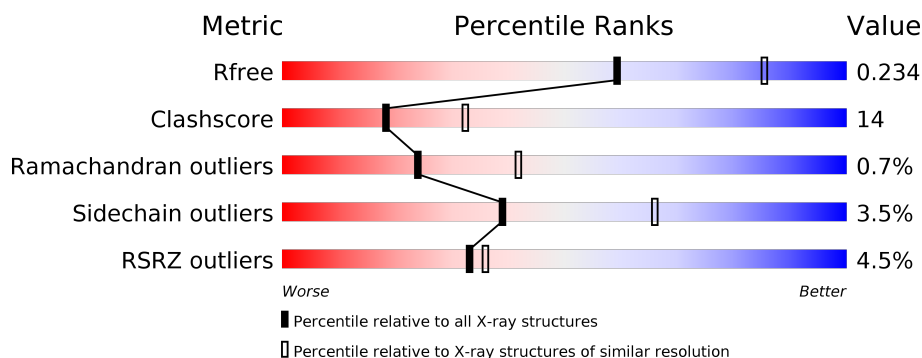
# 1 Overall quality at a glance

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))
$R_{free}$	111664	4155 (2.50-2.50)
Clashscore	122126	4827 (2.50-2.50)
Ramachandran outliers	120053	4735 (2.50-2.50)
Sidechain outliers	120020	4737 (2.50-2.50)
RSRZ outliers	108989	4058 (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 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	25	<div> <div>8%</div> <div> <div></div> <div>20%</div> <div>8%</div> <div>68%</div> </div> </div>
1	B	25	<div> <div>36%</div> <div>56%</div> </div>
2	C	466	<div> <div>4%</div> <div>64%</div> <div>26%</div> <div>8%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3710 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 Large T antigen.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	8	Total	C	N	O	0	0	0
			71	46	15	10			
1	B	11	Total	C	N	O	0	0	0
			82	53	18	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	111	GLY	SER	ENGINEERED	UNP P03070
A	117	ALA	THR	ENGINEERED	UNP P03070
A	120	ALA	SER	ENGINEERED	UNP P03070
A	123	ALA	SER	ENGINEERED	UNP P03070
A	124	ALA	THR	ENGINEERED	UNP P03070
A	134	TYR	ASP	ENGINEERED	UNP P03070
B	111	GLY	SER	ENGINEERED	UNP P03070
B	117	ALA	THR	ENGINEERED	UNP P03070
B	120	ALA	SER	ENGINEERED	UNP P03070
B	123	ALA	SER	ENGINEERED	UNP P03070
B	124	ALA	THR	ENGINEERED	UNP P03070
B	134	TYR	ASP	ENGINEERED	UNP P03070

- Molecule 2 is a protein called Importin alpha-2 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	427	Total	C	N	O	S	0	0	0
			3255	2072	553	620	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	64	MET	-	CLONING ARTIFACT	UNP P03070
C	65	ALA	-	CLONING ARTIFACT	UNP P03070

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Chain	Residue	Modelled	Actual	Comment	Reference
C	66	ASP	-	CLONING ARTIFACT	UNP P03070
C	67	ILE	-	CLONING ARTIFACT	UNP P03070
C	68	GLY	-	CLONING ARTIFACT	UNP P03070
C	69	SER	-	CLONING ARTIFACT	UNP P03070

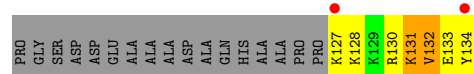
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	6	Total O 6 6	0	0
3	B	16	Total O 16 16	0	0
3	C	280	Total O 280 280	0	0

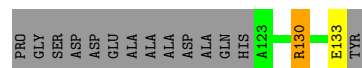
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

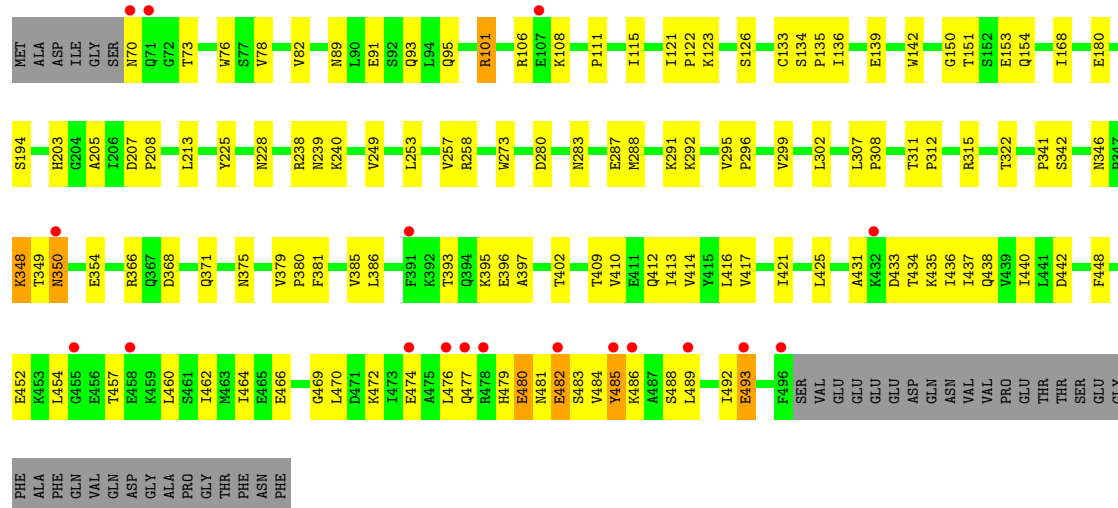
#### • Molecule 1: Large T antigen



#### • Molecule 1: Large T antigen



#### • Molecule 2: Importin alpha-2 subunit



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.16Å 89.28Å 100.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.62 – 2.50 29.62 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (29.62-2.50) 98.3 (29.62-2.50)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.51 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.202 , 0.233 0.203 , 0.234	Depositor DCC
$R_{free}$ test set	1706 reflections (6.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.4	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 47.1	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	0.94	EDS
Total number of atoms	3710	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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	0.53	0/71	2.95	3/88 (3.4%)
1	B	0.46	0/83	0.98	1/108 (0.9%)
2	C	0.33	0/3313	0.58	1/4515 (0.0%)
All	All	0.34	0/3467	0.71	5/4711 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	LYS	N-CA-C	16.49	155.53	111.00
1	A	127	LYS	CB-CA-C	-16.20	78.01	110.40
1	A	128	LYS	N-CA-CB	-11.45	89.99	110.60
1	B	133	GLU	N-CA-C	7.15	130.30	111.00
2	C	348	LYS	CD-CE-NZ	-5.25	99.62	111.70

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	71	0	83	11	0
1	B	82	0	98	1	0
2	C	3255	0	3329	92	0
3	A	6	0	0	1	0
3	B	16	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	280	0	0	11	0
All	All	3710	0	3510	99	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 (99) 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)
2:C:477:GLN:HE21	2:C:489:LEU:HA	1.03	1.15
1:A:131:LYS:O	1:A:133:GLU:N	1.90	1.02
2:C:477:GLN:NE2	2:C:489:LEU:HA	1.84	0.93
2:C:153:GLU:HG2	3:C:801:HOH:O	1.78	0.83
2:C:101:ARG:HD3	2:C:139:GLU:OE1	1.81	0.80
2:C:91:GLU:O	2:C:95:GLN:HG2	1.83	0.79
1:A:132:VAL:O	1:A:132:VAL:HG12	1.88	0.74
2:C:292:LYS:HE2	3:C:562:HOH:O	1.87	0.73
2:C:452:GLU:HG3	2:C:457:THR:HG21	1.69	0.72
2:C:409:THR:H	2:C:412:GLN:NE2	1.93	0.66
2:C:434:THR:O	2:C:438:GLN:HG3	1.94	0.66
2:C:482:GLU:O	2:C:486:LYS:HB2	1.96	0.65
2:C:366:ARG:HD2	2:C:368:ASP:OD1	1.97	0.64
2:C:477:GLN:HE21	2:C:489:LEU:CA	1.96	0.62
1:A:130:ARG:NH2	2:C:396:GLU:OE1	2.32	0.62
2:C:433:ASP:HB3	2:C:436:ILE:HG22	1.82	0.61
2:C:213:LEU:O	2:C:258:ARG:NH2	2.33	0.61
2:C:477:GLN:O	2:C:485:TYR:HB2	2.00	0.61
2:C:477:GLN:HA	2:C:488:SER:HB2	1.82	0.60
2:C:354:GLU:HG3	3:C:742:HOH:O	2.02	0.60
2:C:207:ASP:HB2	2:C:208:PRO:HD3	1.83	0.60
2:C:101:ARG:HD2	2:C:142:TRP:CE3	2.37	0.59
2:C:452:GLU:HA	2:C:457:THR:HG23	1.84	0.58
2:C:472:LYS:O	2:C:476:LEU:HG	2.04	0.57
2:C:299:VAL:O	2:C:302:LEU:HB3	2.05	0.56
1:A:130:ARG:HG2	1:A:130:ARG:HH11	1.70	0.56
2:C:371:GLN:HE22	2:C:375:ASN:HD21	1.54	0.56
2:C:78:VAL:O	2:C:82:VAL:HG23	2.07	0.54
2:C:466:GLU:HB2	3:C:693:HOH:O	2.08	0.54
2:C:311:THR:OG1	3:C:642:HOH:O	2.18	0.54
1:B:130:ARG:NH2	2:C:106:ARG:O	2.41	0.53
2:C:371:GLN:HG2	2:C:412:GLN:OE1	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:379:VAL:HB	2:C:380:PRO:HD3	1.92	0.52
2:C:371:GLN:NE2	2:C:375:ASN:HD21	2.08	0.52
2:C:416:LEU:HG	2:C:421:ILE:HG21	1.93	0.51
2:C:73:THR:HA	2:C:76:TRP:NE1	2.26	0.51
2:C:151:THR:OG1	2:C:154:GLN:HG3	2.11	0.51
2:C:474:GLU:HG2	2:C:492:ILE:HD12	1.93	0.51
2:C:238:ARG:O	2:C:239:ASN:HB2	2.10	0.51
2:C:476:LEU:O	2:C:479:HIS:HB2	2.11	0.50
2:C:462:ILE:O	2:C:466:GLU:HG2	2.11	0.50
2:C:410:VAL:O	2:C:414:VAL:HG23	2.12	0.49
2:C:493:GLU:HG2	3:C:597:HOH:O	2.12	0.49
2:C:438:GLN:O	2:C:442:ASP:HB2	2.12	0.49
2:C:180:GLU:HB2	2:C:225:TYR:CD1	2.48	0.49
2:C:273:TRP:CE2	2:C:312:PRO:HB3	2.48	0.48
2:C:386:LEU:HD22	2:C:421:ILE:HD11	1.95	0.48
2:C:409:THR:H	2:C:412:GLN:HE21	1.60	0.48
2:C:433:ASP:HB3	2:C:436:ILE:CG2	2.44	0.48
2:C:249:VAL:HB	2:C:288:MET:HE1	1.94	0.48
2:C:395:LYS:HE2	3:C:760:HOH:O	2.13	0.48
2:C:205:ALA:O	2:C:208:PRO:HD2	2.14	0.48
2:C:307:LEU:N	2:C:308:PRO:CD	2.77	0.47
1:A:132:VAL:HG22	1:A:134:TYR:HE2	1.79	0.47
2:C:425:LEU:HG	2:C:440:ILE:HG23	1.96	0.47
2:C:287:GLU:O	2:C:291:LYS:HB2	2.14	0.47
2:C:121:ILE:HB	2:C:122:PRO:HD3	1.96	0.47
2:C:194:SER:HB2	3:C:547:HOH:O	2.15	0.47
2:C:350:ASN:ND2	3:C:775:HOH:O	2.48	0.46
2:C:385:VAL:HG13	2:C:393:THR:HG22	1.97	0.46
2:C:452:GLU:C	2:C:454:LEU:H	2.20	0.46
2:C:295:VAL:HB	2:C:296:PRO:HD3	1.97	0.46
2:C:342:SER:HB2	3:C:592:HOH:O	2.15	0.45
2:C:366:ARG:HD3	3:C:728:HOH:O	2.16	0.45
2:C:452:GLU:HA	2:C:457:THR:CG2	2.47	0.45
2:C:240:LYS:HD3	2:C:280:ASP:O	2.17	0.45
2:C:480:GLU:C	2:C:480:GLU:OE1	2.56	0.44
1:A:134:TYR:OH	2:C:315:ARG:NE	2.51	0.44
2:C:95:GLN:H	2:C:95:GLN:HG2	1.62	0.44
2:C:111:PRO:O	2:C:115:ILE:HG12	2.17	0.44
2:C:479:HIS:HD2	2:C:481:ASN:HB3	1.83	0.44
2:C:435:LYS:HD2	2:C:435:LYS:HA	1.76	0.43
1:A:130:ARG:CG	1:A:130:ARG:HH11	2.29	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:VAL:HG23	3:A:138:HOH:O	2.18	0.43
2:C:379:VAL:N	2:C:380:PRO:CD	2.81	0.43
1:A:132:VAL:O	1:A:132:VAL:CG1	2.60	0.43
2:C:413:ILE:O	2:C:417:VAL:HG23	2.18	0.43
2:C:431:ALA:O	2:C:437:ILE:HD11	2.18	0.43
2:C:346:ASN:HD22	2:C:348:LYS:H	1.67	0.43
2:C:150:GLY:HA3	2:C:154:GLN:OE1	2.19	0.42
2:C:134:SER:N	2:C:135:PRO:CD	2.82	0.42
2:C:168:ILE:HD12	2:C:203:HIS:HB3	2.01	0.42
2:C:89:ASN:O	2:C:93:GLN:HG3	2.20	0.42
1:A:131:LYS:HA	2:C:322:THR:HG21	2.01	0.42
2:C:452:GLU:HG3	2:C:457:THR:CG2	2.45	0.42
2:C:477:GLN:HA	2:C:488:SER:CB	2.48	0.42
2:C:341:PRO:HB3	2:C:381:PHE:CZ	2.55	0.41
2:C:470:LEU:C	2:C:472:LYS:H	2.23	0.41
2:C:283:ASN:N	2:C:283:ASN:HD22	2.18	0.41
2:C:433:ASP:O	2:C:436:ILE:HG22	2.20	0.41
2:C:436:ILE:O	2:C:440:ILE:HG13	2.21	0.41
2:C:464:ILE:O	2:C:469:GLY:N	2.44	0.41
2:C:123:LYS:O	2:C:126:SER:HB3	2.21	0.41
2:C:133:CYS:O	2:C:136:ILE:HG22	2.20	0.41
2:C:385:VAL:HG12	2:C:397:ALA:HB2	2.02	0.41
1:A:134:TYR:OH	2:C:315:ARG:CZ	2.69	0.41
2:C:348:LYS:N	2:C:348:LYS:HD2	2.36	0.41
2:C:253:LEU:O	2:C:257:VAL:HG23	2.22	0.40
2:C:448:PHE:CD1	2:C:460:LEU:HD23	2.57	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	6/25 (24%)	3 (50%)	1 (17%)	2 (33%)	0	0
1	B	9/25 (36%)	9 (100%)	0	0	100	100
2	C	425/466 (91%)	410 (96%)	14 (3%)	1 (0%)	49	71
All	All	440/516 (85%)	422 (96%)	15 (3%)	3 (1%)	24	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	VAL
1	A	131	LYS
2	C	484	VAL

### 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	7/18 (39%)	7 (100%)	0	100	100
1	B	8/18 (44%)	7 (88%)	1 (12%)	5	9
2	C	358/390 (92%)	346 (97%)	12 (3%)	40	67
All	All	373/426 (88%)	360 (96%)	13 (4%)	39	66

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	130	ARG
2	C	70	ASN
2	C	101	ARG
2	C	108	LYS
2	C	228	ASN
2	C	349	THR
2	C	350	ASN
2	C	402	THR
2	C	480	GLU
2	C	482	GLU

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Mol	Chain	Res	Type
2	C	483	SER
2	C	485	TYR
2	C	493	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	86	ASN
2	C	283	ASN
2	C	346	ASN
2	C	371	GLN
2	C	375	ASN
2	C	412	GLN
2	C	438	GLN
2	C	446	ASN
2	C	477	GLN
2	C	479	HIS

### 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 ⓘ

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	8/25 (32%)	1.81	2 (25%) 0 0	51, 64, 77, 83	0
1	B	11/25 (44%)	-0.14	0 100 100	24, 35, 48, 59	0
2	C	427/466 (91%)	-0.00	18 (4%) 36 39	18, 32, 76, 94	0
All	All	446/516 (86%)	0.03	20 (4%) 33 36	18, 33, 77, 94	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	485	TYR	6.6
2	C	71	GLN	6.4
1	A	127	LYS	5.3
1	A	134	TYR	4.9
2	C	489	LEU	4.4
2	C	496	PHE	4.4
2	C	482	GLU	3.9
2	C	474	GLU	3.7
2	C	493	GLU	3.6
2	C	455	GLY	3.3
2	C	70	ASN	3.3
2	C	432	LYS	3.2
2	C	478	ARG	2.9
2	C	458	GLU	2.7
2	C	350	ASN	2.6
2	C	476	LEU	2.2
2	C	391	PHE	2.2
2	C	477	GLN	2.2
2	C	486	LYS	2.2
2	C	107	GLU	2.2

## 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.