



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:03 AM GMT

PDB ID : 2VS6
Title : K173A, R174A, K177A-TRICHOSANTHIN
Authors : Too, P.H.; Ma, M.K.; Mak, A.N.; Tung, C.K.; Zhu, G.; Au, S.W.; Wong, K.B.;
Shaw, P.C.; Ng, A.
Deposited on : 2008-04-21
Resolution : 2.40 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

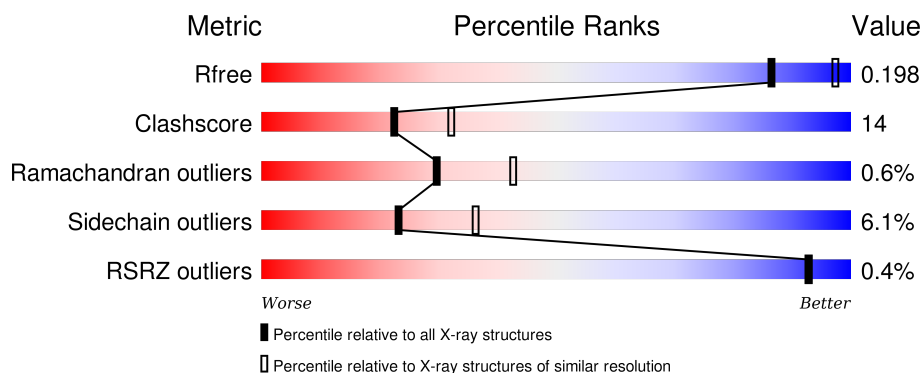
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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 ≥ 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	248	 74% 21% •
1	B	248	 76% 21% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4184 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 RIBOSOME-INACTIVATING PROTEIN ALPHA-TRICHOSANTHIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1896	1202	320	369	5			
1	B	248	Total	C	N	O	S	0	0	0
			1907	1208	324	370	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	LYS	ENGINEERED MUTATION	UNP P09989
A	174	ALA	ARG	ENGINEERED MUTATION	UNP P09989
A	177	ALA	LYS	ENGINEERED MUTATION	UNP P09989
B	173	ALA	LYS	ENGINEERED MUTATION	UNP P09989
B	174	ALA	ARG	ENGINEERED MUTATION	UNP P09989
B	177	ALA	LYS	ENGINEERED MUTATION	UNP P09989

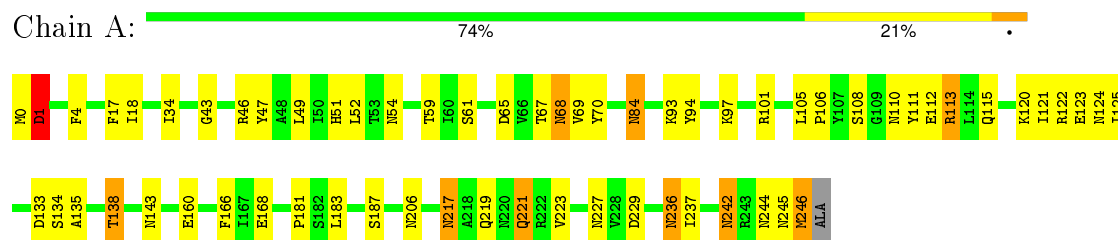
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	215	Total	O	0	0
			215	215		
2	B	166	Total	O	0	0
			166	166		

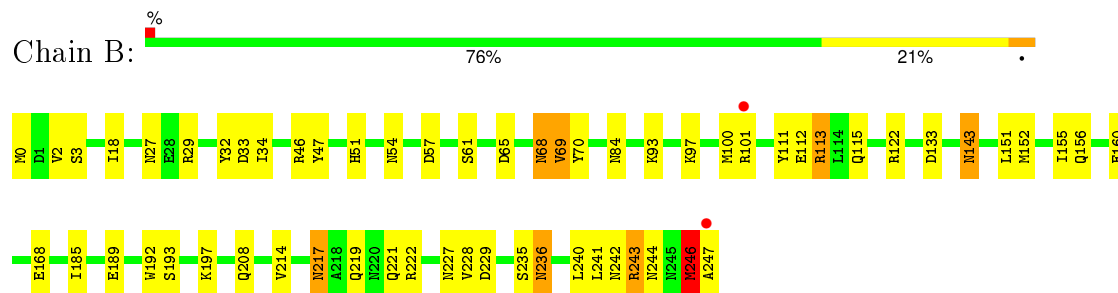
3 Residue-property plots

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: RIBOSOME-INACTIVATING PROTEIN ALPHA-TRICHOSANTHIN



• Molecule 1: RIBOSOME-INACTIVATING PROTEIN ALPHA-TRICHOSANTHIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.03Å 95.84Å 101.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 30.54 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (50.00-2.40) 98.8 (30.54-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.191 , 0.236 0.195 , 0.198	Depositor DCC
R_{free} test set	1354 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 54252 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4184	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	9.22	5/1928 (0.3%)	1.22	5/2623 (0.2%)
1	B	0.66	7/1939 (0.4%)	0.68	4/2637 (0.2%)
All	All	6.53	12/3867 (0.3%)	0.99	9/5260 (0.2%)

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	2	1
1	B	1	0
All	All	3	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	246	MET	C-O	351.49	7.91	1.23
1	A	246	MET	CA-C	200.03	6.73	1.52
1	B	246	MET	C-O	-13.86	0.97	1.23
1	B	0	MET	C-O	-8.40	1.07	1.23
1	B	0	MET	N-CA	-8.17	1.30	1.46
1	A	0	MET	CA-CB	-7.56	1.37	1.53
1	B	246	MET	CB-CG	-6.78	1.29	1.51
1	A	0	MET	C-O	-6.75	1.10	1.23
1	B	246	MET	N-CA	-6.52	1.33	1.46
1	B	246	MET	CA-C	-6.22	1.36	1.52
1	B	0	MET	CA-C	-6.13	1.37	1.52
1	A	0	MET	CB-CG	-5.47	1.33	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	MET	CA-C-O	-50.85	13.32	120.10
1	A	246	MET	N-CA-C	-13.62	74.23	111.00
1	B	0	MET	CG-SD-CE	9.59	115.55	100.20
1	B	246	MET	CB-CG-SD	-7.30	90.49	112.40
1	A	1	ASP	O-C-N	-6.41	112.44	122.70
1	A	246	MET	N-CA-CB	5.83	121.09	110.60
1	B	0	MET	CB-CA-C	5.75	121.89	110.40
1	B	0	MET	CB-CG-SD	5.22	128.06	112.40
1	A	1	ASP	N-CA-CB	5.05	119.68	110.60

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	1	ASP	CA
1	A	246	MET	CA
1	B	246	MET	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	ASP	Mainchain

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	1896	0	1901	59	0
1	B	1907	0	1917	45	0
2	A	215	0	0	0	0
2	B	166	0	0	2	0
All	All	4184	0	3818	104	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 (104) 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:68:ASN:HB3	1:A:246:MET:C	1.27	1.49
1:A:245:ASN:ND2	1:A:246:MET:C	1.84	1.28
1:A:246:MET:HA	1:A:246:MET:O	1.39	1.22
1:A:68:ASN:CB	1:A:246:MET:C	2.12	1.15
1:A:120:LYS:HD3	1:A:125:ILE:HD11	1.37	1.07
1:A:246:MET:O	1:A:246:MET:CA	2.06	1.04
1:A:245:ASN:HD21	1:A:246:MET:C	1.59	1.00
1:A:242:ASN:HD22	1:A:244:ASN:H	1.11	0.96
1:A:227:ASN:HD22	1:A:229:ASP:H	1.11	0.94
1:A:245:ASN:HD22	1:A:246:MET:C	1.70	0.91
1:B:242:ASN:HD22	1:B:244:ASN:H	1.21	0.88
1:B:227:ASN:HD22	1:B:229:ASP:H	1.21	0.88
1:B:246:MET:O	1:B:246:MET:HG2	1.73	0.88
1:B:115:GLN:NE2	1:B:122:ARG:H	1.74	0.86
1:A:120:LYS:HD3	1:A:125:ILE:CD1	2.10	0.81
1:B:115:GLN:HE21	1:B:122:ARG:H	1.29	0.80
1:B:217:ASN:ND2	1:B:221:GLN:H	1.80	0.79
1:B:65:ASP:HB3	1:B:68:ASN:HD21	1.46	0.78
1:A:217:ASN:HD21	1:A:221:GLN:HG3	1.55	0.71
1:A:115:GLN:HE21	1:A:122:ARG:H	1.38	0.71
1:A:111:TYR:O	1:A:115:GLN:HG3	1.90	0.71
1:B:227:ASN:ND2	1:B:229:ASP:H	1.89	0.70
1:A:227:ASN:ND2	1:A:229:ASP:HB2	2.08	0.69
1:B:243:ARG:HA	1:B:247:ALA:HA	1.74	0.69
1:A:115:GLN:NE2	1:A:122:ARG:H	1.90	0.68
1:B:93:LYS:O	1:B:97:LYS:HE3	1.94	0.68
1:B:217:ASN:HD21	1:B:221:GLN:H	1.40	0.68
1:A:134:SER:O	1:A:138:THR:HG22	1.97	0.64
1:B:68:ASN:ND2	1:B:70:TYR:H	1.98	0.62
1:A:166:PHE:H	1:A:236:ASN:ND2	1.99	0.61
1:A:227:ASN:ND2	1:A:229:ASP:H	1.91	0.60
1:A:242:ASN:ND2	1:A:244:ASN:H	1.92	0.60
1:A:227:ASN:HD21	1:A:229:ASP:HB2	1.65	0.60
1:A:65:ASP:HB3	1:A:68:ASN:HD21	1.65	0.60
1:B:197:LYS:HD2	1:B:247:ALA:HB1	1.82	0.60
1:A:217:ASN:C	1:A:217:ASN:HD22	2.03	0.59
1:A:217:ASN:HD21	1:A:221:GLN:H	1.52	0.58
1:B:54:ASN:HB2	1:B:133:ASP:OD1	2.04	0.57
1:B:246:MET:CG	1:B:246:MET:O	2.29	0.57
1:B:217:ASN:C	1:B:217:ASN:HD22	2.09	0.56
1:B:65:ASP:HB3	1:B:68:ASN:ND2	2.20	0.56
1:B:112:GLU:OE1	1:B:113:ARG:NE	2.38	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ILE:HG13	1:B:168:GLU:HG3	1.87	0.56
1:A:166:PHE:H	1:A:236:ASN:HD21	1.53	0.55
1:A:217:ASN:ND2	1:A:221:GLN:H	2.06	0.54
1:A:108:SER:OG	1:A:113:ARG:HG2	2.06	0.53
1:A:68:ASN:HD22	1:A:68:ASN:C	2.12	0.52
1:A:110:ASN:ND2	1:A:112:GLU:HB3	2.24	0.52
1:B:100:MET:O	1:B:101:ARG:HD2	2.10	0.51
1:A:68:ASN:ND2	1:A:70:TYR:H	2.08	0.50
1:A:1:ASP:OD1	1:A:49:LEU:N	2.38	0.50
1:A:54:ASN:HB2	1:A:133:ASP:OD1	2.14	0.48
1:A:242:ASN:HD22	1:A:244:ASN:N	1.95	0.48
1:B:217:ASN:ND2	1:B:219:GLN:H	2.11	0.48
1:A:68:ASN:HD22	1:A:69:VAL:N	2.12	0.48
1:A:227:ASN:HD22	1:A:229:ASP:N	1.94	0.48
1:A:217:ASN:ND2	1:A:221:GLN:HG3	2.28	0.48
1:A:123:GLU:HA	1:A:181:PRO:HG2	1.96	0.47
1:B:227:ASN:ND2	1:B:229:ASP:HB2	2.29	0.47
1:B:68:ASN:C	1:B:68:ASN:HD22	2.16	0.47
1:B:214:VAL:HG11	1:B:222:ARG:NH2	2.29	0.47
1:B:97:LYS:NZ	2:B:2071:HOH:O	2.47	0.47
1:B:68:ASN:HD22	1:B:69:VAL:N	2.13	0.47
1:A:51:HIS:HD2	1:A:61:SER:OG	1.98	0.47
1:A:120:LYS:CD	1:A:125:ILE:HD11	2.27	0.46
1:A:93:LYS:O	1:A:97:LYS:HD3	2.15	0.46
1:B:192:TRP:CH2	1:B:241:LEU:HB2	2.51	0.46
1:B:32:TYR:CE2	1:B:228:VAL:HG11	2.50	0.46
1:B:242:ASN:ND2	1:B:244:ASN:H	2.01	0.46
1:B:143:ASN:ND2	2:B:2106:HOH:O	2.49	0.45
1:B:27:ASN:HD22	1:B:27:ASN:N	2.13	0.45
1:A:68:ASN:CA	1:A:246:MET:C	2.81	0.45
1:B:185:ILE:O	1:B:189:GLU:HG3	2.16	0.45
1:A:217:ASN:C	1:A:217:ASN:ND2	2.70	0.45
1:B:51:HIS:HD2	1:B:61:SER:OG	2.00	0.45
1:B:160:GLU:OE1	1:B:160:GLU:HA	2.17	0.45
1:B:151:LEU:O	1:B:155:ILE:HG13	2.16	0.45
1:A:43:GLY:HA3	1:A:94:TYR:CZ	2.53	0.44
1:A:237:ILE:O	1:A:237:ILE:HG23	2.17	0.44
1:A:67:THR:C	1:A:246:MET:C	2.76	0.44
1:A:4:PHE:HB2	1:A:17:PHE:CG	2.53	0.44
1:B:46:ARG:HD2	1:B:47:TYR:CZ	2.53	0.44
1:B:2:VAL:CG1	1:B:3:SER:N	2.81	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:ASN:HD22	1:B:236:ASN:C	2.21	0.44
1:B:246:MET:O	1:B:247:ALA:HB2	2.18	0.43
1:B:111:TYR:O	1:B:115:GLN:HG3	2.18	0.43
1:A:68:ASN:CG	1:A:246:MET:C	2.76	0.43
1:A:217:ASN:ND2	1:A:219:GLN:H	2.16	0.43
1:A:113:ARG:N	1:A:113:ARG:HD2	2.34	0.43
1:A:18:ILE:HG13	1:A:168:GLU:HG3	2.01	0.43
1:A:121:ILE:HG13	1:A:123:GLU:HG2	2.01	0.42
1:A:84:ASN:HB2	1:A:105:LEU:O	2.19	0.42
1:B:222:ARG:HB3	1:B:222:ARG:NH1	2.34	0.42
1:A:34:ILE:HD13	1:A:237:ILE:HG23	2.01	0.42
1:A:183:LEU:O	1:A:187:SER:HB2	2.18	0.42
1:B:34:ILE:HG22	1:B:240:LEU:HG	2.01	0.42
1:A:46:ARG:HD2	1:A:47:TYR:CZ	2.55	0.41
1:B:152:MET:O	1:B:156:GLN:HG3	2.20	0.41
1:A:135:ALA:O	1:A:138:THR:HG23	2.21	0.40
1:B:32:TYR:O	1:B:33:ASP:HB2	2.21	0.40
1:A:52:LEU:O	1:A:59:THR:HG23	2.22	0.40
1:B:68:ASN:HD22	1:B:70:TYR:H	1.67	0.40
1:A:105:LEU:HA	1:A:106:PRO:HD3	1.89	0.40
1:A:160:GLU:OE1	1:A:160:GLU:HA	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	245/248 (99%)	235 (96%)	10 (4%)	0	100	100
1	B	246/248 (99%)	238 (97%)	5 (2%)	3 (1%)	16	23
All	All	491/496 (99%)	473 (96%)	15 (3%)	3 (1%)	30	43

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	246	MET
1	B	29	ARG
1	B	69	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	205/206 (100%)	192 (94%)	13 (6%)	22	35
1	B	206/206 (100%)	194 (94%)	12 (6%)	25	39
All	All	411/412 (100%)	386 (94%)	25 (6%)	23	36

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	84	ASN
1	A	101	ARG
1	A	113	ARG
1	A	124	ASN
1	A	138	THR
1	A	143	ASN
1	A	206	ASN
1	A	217	ASN
1	A	221	GLN
1	A	223	VAL
1	A	236	ASN
1	A	242	ASN
1	B	57	ASP
1	B	68	ASN
1	B	84	ASN
1	B	113	ARG
1	B	143	ASN
1	B	193	SER
1	B	208	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	217	ASN
1	B	235	SER
1	B	236	ASN
1	B	243	ARG
1	B	246	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	51	HIS
1	A	68	ASN
1	A	84	ASN
1	A	115	GLN
1	A	124	ASN
1	A	143	ASN
1	A	145	ASN
1	A	156	GLN
1	A	206	ASN
1	A	217	ASN
1	A	220	ASN
1	A	221	GLN
1	A	227	ASN
1	A	236	ASN
1	A	242	ASN
1	A	244	ASN
1	A	245	ASN
1	B	20	ASN
1	B	27	ASN
1	B	51	HIS
1	B	68	ASN
1	B	84	ASN
1	B	110	ASN
1	B	115	GLN
1	B	124	ASN
1	B	143	ASN
1	B	145	ASN
1	B	156	GLN
1	B	170	GLN
1	B	217	ASN
1	B	221	GLN
1	B	227	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	236	ASN
1	B	242	ASN
1	B	244	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, 95th 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 ²)	Q<0.9
1	A	247/248 (99%)	-0.75	0 100 100	9, 18, 30, 46	0
1	B	248/248 (100%)	-0.64	2 (0%) 87 87	9, 18, 37, 55	0
All	All	495/496 (99%)	-0.69	2 (0%) 93 93	9, 18, 34, 55	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	247	ALA	3.6
1	B	101	ARG	2.9

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.