



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4WKR / pdb_00004wkr
Title : LaRP7 wrapping up the 3' hairpin of 7SK non-coding RNA (302-332)
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Deposited on : 2014-10-03
Resolution : 3.20 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

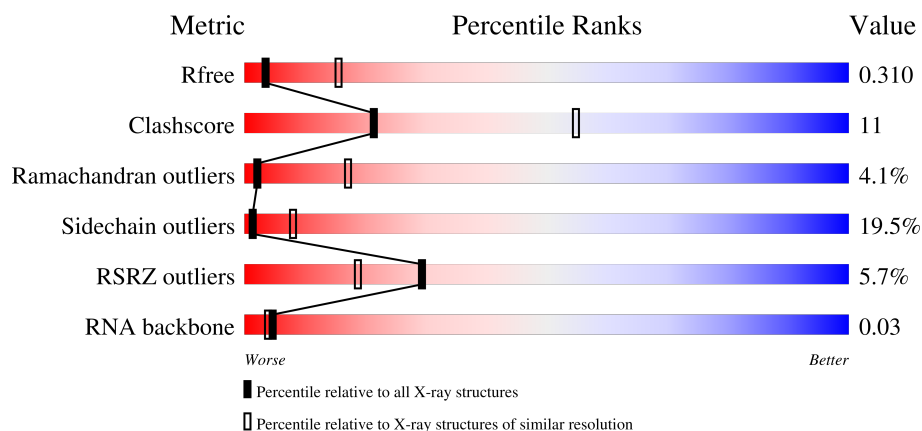
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)
RNA backbone	3983	1222 (3.50-2.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	208	3% 48% 21% 7% • 23%
1	B	208	5% 46% 23% 8% • 23%
2	C	31	6% 10% 84%
2	D	31	3% • • 10% 84%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2701 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 La-related protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1261	803	224	232	2			
1	B	161	Total	C	N	O	S	0	0	0
			1267	808	227	230	2			

- Molecule 2 is a RNA chain called 7SK GGHP4 (300-332).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	5	Total	C	N	O	P	0	0	0
			68	27	6	30	5			
2	D	5	Total	C	N	O	P	0	0	0
			100	45	11	39	5			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	O	0	0
			3	3		
3	C	1	Total	O	0	0
			1	1		
3	D	1	Total	O	0	0
			1	1		



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.45Å 33.50Å 119.08Å 90.00° 128.99° 90.00°	Depositor
Resolution (Å)	59.06 – 3.20 59.06 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.0 (59.06-3.20) 94.0 (59.06-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.221 , 0.274 0.267 , 0.310	Depositor DCC
R_{free} test set	382 reflections (4.40%)	wwPDB-VP
Wilson B-factor (Å ²)	90.5	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 68.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2701	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.31% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.99	1/1286 (0.1%)	1.99	32/1741 (1.8%)
1	B	1.24	4/1291 (0.3%)	1.76	23/1743 (1.3%)
2	C	0.90	0/71	1.29	0/104
2	D	1.36	2/109 (1.8%)	1.83	6/166 (3.6%)
All	All	1.13	7/2757 (0.3%)	1.86	61/3754 (1.6%)

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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	188	ASN	C-N	24.63	1.74	1.34
1	B	104	LEU	C-N	-16.73	1.08	1.33
1	A	68	TYR	C-N	-16.18	1.13	1.33
1	B	105	GLU	C-N	11.07	1.47	1.33
2	D	-5	C	C1'-N1	9.23	1.62	1.48
1	B	105	GLU	CA-C	5.75	1.60	1.52
2	D	-4	U	C1'-N1	5.34	1.55	1.47

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	105	GLU	N-CA-C	24.75	143.13	113.18
1	A	104	LEU	CA-C-N	24.71	168.74	121.54
1	A	104	LEU	C-N-CA	24.71	168.74	121.54
1	A	68	TYR	N-CA-C	19.06	138.20	111.52
1	A	105	GLU	N-CA-CB	-16.39	82.79	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ASN	O-C-N	-15.38	99.28	121.12
1	B	104	LEU	O-C-N	-14.56	104.18	122.09
1	A	68	TYR	O-C-N	-11.76	108.47	122.58
1	B	102	LEU	N-CA-C	-10.80	92.49	109.23
1	B	105	GLU	CB-CA-C	-10.72	88.62	109.95
1	B	161	SER	N-CA-C	-10.40	88.64	110.80
1	A	161	SER	N-CA-C	-10.26	88.94	110.80
1	A	68	TYR	CB-CA-C	-9.59	99.01	111.86
1	A	105	GLU	N-CA-C	8.95	129.87	110.80
1	A	69	VAL	N-CA-C	8.90	120.96	109.30
1	A	161	SER	CB-CA-C	8.82	127.97	110.42
1	B	38	ILE	N-CA-C	-8.77	102.18	110.42
2	D	-4	U	P-O3'-C3'	8.75	133.32	120.20
1	A	102	LEU	N-CA-C	8.60	123.87	113.12
1	B	161	SER	CB-CA-C	8.56	127.45	110.42
1	A	104	LEU	CB-CA-C	-8.14	98.44	110.16
1	B	105	GLU	CA-C-O	7.67	128.93	119.11
1	A	68	TYR	CA-C-N	7.24	131.25	120.69
1	A	68	TYR	C-N-CA	7.24	131.25	120.69
1	A	38	ILE	N-CA-C	-6.98	104.05	110.82
1	B	43	ASP	CA-CB-CG	6.76	119.36	112.60
1	B	105	GLU	O-C-N	6.57	132.02	122.43
1	A	162	THR	N-CA-CB	-6.44	99.61	110.49
1	B	162	THR	N-CA-CB	-6.39	99.68	110.49
2	D	-5	C	C3'-C2'-C1'	-6.16	95.14	101.30
1	A	112	LYS	CB-CA-C	-6.12	101.25	110.06
1	B	160	LYS	N-CA-C	-6.05	102.97	110.41
1	B	101	GLU	N-CA-CB	5.96	120.61	111.56
1	A	43	ASP	CA-CB-CG	5.91	118.51	112.60
1	A	160	LYS	N-CA-C	-5.82	103.25	110.41
1	A	146	GLY	CA-C-N	5.82	130.00	120.63
1	A	146	GLY	C-N-CA	5.82	130.00	120.63
1	A	67	GLY	N-CA-C	-5.78	102.16	112.22
1	B	146	GLY	CA-C-N	5.74	129.87	120.63
1	B	146	GLY	C-N-CA	5.74	129.87	120.63
1	B	109	ILE	N-CA-CB	5.45	118.94	111.41
1	B	71	ILE	CA-C-N	5.41	127.47	120.44
1	B	71	ILE	C-N-CA	5.41	127.47	120.44
2	D	-5	C	C2'-C3'-O3'	5.39	121.79	113.70
1	A	143	ARG	CA-C-N	5.37	127.81	120.46
1	A	143	ARG	C-N-CA	5.37	127.81	120.46
1	B	162	THR	N-CA-C	5.33	122.15	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	LYS	CA-C-N	5.32	131.69	121.54
1	A	63	LYS	C-N-CA	5.32	131.69	121.54
1	B	112	LYS	CB-CA-C	-5.31	99.86	110.42
2	D	-5	C	C4'-C3'-O3'	5.31	120.96	113.00
1	B	99	VAL	N-CA-C	5.22	113.89	106.53
1	B	131	LEU	N-CA-C	5.20	118.48	110.42
1	A	162	THR	N-CA-C	5.18	121.84	110.80
1	B	188	ASN	O-C-N	5.18	127.24	121.43
2	D	-4	U	C3'-C2'-C1'	5.18	106.68	101.50
2	D	-2	U	P-O5'-C5'	5.11	128.57	120.90
1	A	71	ILE	CA-C-N	5.09	127.06	120.44
1	A	71	ILE	C-N-CA	5.09	127.06	120.44
1	A	142	GLU	CB-CG-CD	5.05	121.19	112.60
1	A	109	ILE	N-CA-CB	5.02	118.34	111.41

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	ASN	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	1261	0	1225	20	0
1	B	1267	0	1249	37	0
2	C	68	0	31	0	0
2	D	100	0	52	0	0
3	A	3	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
All	All	2701	0	2557	57	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 11.

All (57) 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:188:ASN:C	1:B:189:PRO:N	1.74	1.45
1:B:105:GLU:HB2	1:B:107:THR:HG22	1.18	1.16
1:B:105:GLU:HB2	1:B:107:THR:CG2	2.05	0.84
1:B:188:ASN:ND2	1:B:189:PRO:CD	2.46	0.79
1:B:105:GLU:CB	1:B:107:THR:HG22	2.07	0.79
1:B:97:SER:OG	1:B:100:VAL:HG12	1.84	0.76
1:B:188:ASN:CA	1:B:189:PRO:N	2.48	0.75
1:B:188:ASN:ND2	1:B:189:PRO:HD2	2.05	0.71
1:B:188:ASN:ND2	1:B:189:PRO:HD3	2.05	0.71
1:A:52:HIS:HB3	1:A:118:ARG:HD2	1.71	0.71
1:B:99:VAL:O	1:B:100:VAL:HB	1.92	0.68
1:B:188:ASN:CG	1:B:189:PRO:HD2	2.20	0.67
1:B:188:ASN:C	1:B:189:PRO:CA	2.68	0.65
1:A:100:VAL:HA	1:A:110:ARG:O	1.96	0.64
1:A:188:ASN:HB3	1:A:189:PRO:CD	2.26	0.64
1:B:100:VAL:O	1:B:112:LYS:HD2	2.00	0.62
1:A:99:VAL:O	1:A:100:VAL:HG12	2.01	0.61
1:B:99:VAL:HG21	1:B:111:ARG:HH21	1.66	0.60
1:B:188:ASN:CG	1:B:189:PRO:CD	2.74	0.60
1:B:99:VAL:O	1:B:99:VAL:HG13	2.03	0.59
1:A:178:GLN:HA	1:A:181:LYS:HE3	1.86	0.58
1:B:175:THR:C	1:B:177:GLU:H	2.13	0.55
1:A:188:ASN:HB3	1:A:189:PRO:HD3	1.88	0.55
1:A:102:LEU:HG	1:A:109:ILE:HG22	1.88	0.54
1:B:102:LEU:HG	1:B:109:ILE:HG22	1.88	0.54
1:A:140:TRP:HD1	1:A:143:ARG:HH21	1.54	0.54
1:A:105:GLU:HB3	1:A:107:THR:HG22	1.89	0.53
1:B:52:HIS:HB3	1:B:118:ARG:HD3	1.91	0.53
1:B:134:LYS:C	1:B:136:VAL:H	2.18	0.52
1:A:134:LYS:C	1:A:136:VAL:H	2.18	0.51
1:A:188:ASN:O	1:A:189:PRO:C	2.55	0.50
1:B:188:ASN:C	1:B:189:PRO:C	2.81	0.49
1:B:52:HIS:HB3	1:B:118:ARG:CD	2.42	0.48
1:B:183:ILE:O	1:B:186:LEU:HB3	2.13	0.48
1:A:183:ILE:O	1:A:186:LEU:HB3	2.13	0.48
1:B:103:ASP:OD1	1:B:103:ASP:N	2.24	0.48
1:B:126:THR:HA	1:B:172:GLU:HA	1.95	0.48
1:A:84:THR:HG22	1:A:86:ASP:H	1.79	0.47
1:A:52:HIS:HB3	1:A:118:ARG:CD	2.44	0.47
1:B:84:THR:HG22	1:B:86:ASP:H	1.79	0.47
1:B:140:TRP:O	1:B:144:VAL:HG22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ASN:HA	1:B:189:PRO:N	2.31	0.46
1:B:126:THR:HG22	1:B:172:GLU:HB2	1.97	0.46
1:B:105:GLU:CB	1:B:107:THR:CG2	2.81	0.45
1:B:155:SER:HB3	1:B:170:PHE:HB2	1.98	0.45
1:A:157:PRO:HB2	1:A:166:LYS:HD2	1.99	0.44
1:A:155:SER:HB3	1:A:170:PHE:HB2	1.99	0.44
1:A:152:VAL:HG21	1:A:174:GLU:HG2	2.00	0.43
1:B:68:TYR:HB3	1:B:108:ARG:HB3	2.01	0.43
1:B:154:ILE:HG12	1:B:171:VAL:HG13	2.01	0.42
1:B:105:GLU:N	1:B:105:GLU:OE1	2.53	0.42
1:A:154:ILE:HG12	1:A:171:VAL:HG13	2.01	0.42
1:A:87:GLY:HA2	1:A:90:ILE:HD12	2.01	0.42
1:B:91:ALA:HB1	1:B:95:ARG:HH12	1.85	0.41
1:B:87:GLY:HA2	1:B:90:ILE:HD12	2.02	0.41
1:A:135:ASN:C	1:A:137:ASN:H	2.28	0.40
1:B:135:ASN:C	1:B:137:ASN:H	2.29	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	159/208 (76%)	139 (87%)	13 (8%)	7 (4%)	2	15
1	B	159/208 (76%)	132 (83%)	21 (13%)	6 (4%)	2	18
All	All	318/416 (76%)	271 (85%)	34 (11%)	13 (4%)	2	17

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	VAL
1	B	136	VAL

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Mol	Chain	Res	Type
1	A	48	ASP
1	A	64	SER
1	B	48	ASP
1	B	100	VAL
1	A	137	ASN
1	B	137	ASN
1	A	100	VAL
1	B	68	TYR
1	A	65	ARG
1	A	135	ASN
1	B	135	ASN

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	130/186 (70%)	106 (82%)	24 (18%)	1	9
1	B	131/186 (70%)	104 (79%)	27 (21%)	1	7
All	All	261/372 (70%)	210 (80%)	51 (20%)	1	8

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	51	LEU
1	A	54	ASP
1	A	61	ILE
1	A	65	ARG
1	A	69	VAL
1	A	73	LEU
1	A	84	THR
1	A	89	LEU
1	A	100	VAL
1	A	103	ASP
1	A	104	LEU

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Mol	Chain	Res	Type
1	A	105	GLU
1	A	109	ILE
1	A	110	ARG
1	A	111	ARG
1	A	113	LYS
1	A	132	LEU
1	A	164	ASP
1	A	171	VAL
1	A	174	GLU
1	A	175	THR
1	A	183	ILE
1	A	186	LEU
1	B	40	LYS
1	B	41	GLN
1	B	43	ASP
1	B	51	LEU
1	B	54	ASP
1	B	60	GLN
1	B	61	ILE
1	B	64	SER
1	B	73	LEU
1	B	79	LYS
1	B	84	THR
1	B	89	LEU
1	B	100	VAL
1	B	103	ASP
1	B	109	ILE
1	B	110	ARG
1	B	111	ARG
1	B	113	LYS
1	B	118	ARG
1	B	132	LEU
1	B	147	LYS
1	B	164	ASP
1	B	171	VAL
1	B	174	GLU
1	B	183	ILE
1	B	184	GLU
1	B	186	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	137	ASN
1	A	187	ASN
1	B	41	GLN
1	B	52	HIS
1	B	187	ASN
1	B	188	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	3/31 (9%)	2 (66%)	1 (33%)
2	D	5/31 (16%)	3 (60%)	3 (60%)
All	All	8/62 (12%)	5 (62%)	4 (50%)

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	-2	U
2	C	-1	U
2	D	-4	U
2	D	-3	U
2	D	-2	U

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	-3	U
2	D	-5	C
2	D	-4	U
2	D	-3	U

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 oligosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	2
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	188:ASN	C	189:PRO	N	1.74
1	A	68:TYR	C	69:VAL	N	1.13
1	B	104:LEU	C	105:GLU	N	1.08

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	161/208 (77%)	0.47	7 (4%) 40 24	45, 86, 123, 139	0
1	B	161/208 (77%)	0.61	11 (6%) 23 15	43, 92, 128, 160	0
2	C	5/31 (16%)	0.39	0 100 100	93, 99, 139, 165	0
2	D	5/31 (16%)	0.96	1 (20%) 3 2	100, 106, 138, 155	0
All	All	332/478 (69%)	0.55	19 (5%) 29 19	43, 90, 129, 165	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	106	GLY	5.4
1	B	145	PHE	5.3
1	B	105	GLU	4.0
1	A	161	SER	3.4
1	A	189	PRO	3.0
1	A	145	PHE	2.9
1	B	171	VAL	2.9
1	A	135	ASN	2.8
1	B	175	THR	2.7
1	A	168	PHE	2.6
1	A	141	ILE	2.6
1	B	141	ILE	2.6
1	A	105	GLU	2.4
1	B	51	LEU	2.3
1	B	170	PHE	2.2
1	B	73	LEU	2.2
2	D	-5	C	2.1
1	B	155	SER	2.1
1	B	161	SER	2.1

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 oligosaccharides 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.