



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:57 PM UTC

PDB ID : 5N94 / pdb_00005n94
Title : Crystal Structure of Drosophila DHX36 helicase in complex with polyU
Authors : Chen, W.-F.; Rety, S.; Hai-Lei Guo, H.-L.; Wu, W.-Q.; Liu, N.-N.; Liu, Q.-W.;
Dai, Y.-X.; Xi, X.-G.
Deposited on : 2017-02-24
Resolution : 2.43 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

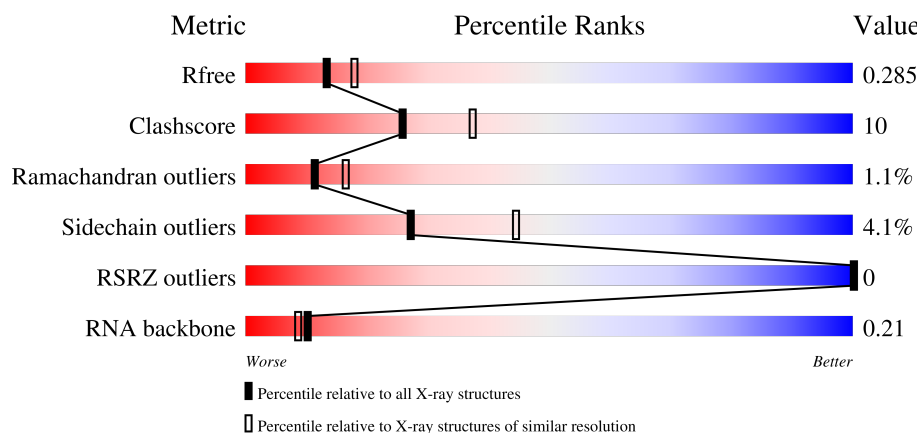
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

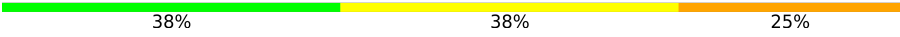
The reported resolution of this entry is 2.43 Å.

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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)
RNA backbone	3983	1058 (2.70-2.14)

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	944	
1	B	944	
2	C	8	
2	D	8	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14015 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 CG9323, isoform A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	849	Total	C	N	O	S	0	0	0
			6814	4304	1200	1265	45			
1	B	849	Total	C	N	O	S	0	0	0
			6814	4304	1200	1265	45			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	943	VAL	-	expression tag	UNP Q8SWT2
A	944	ASP	-	expression tag	UNP Q8SWT2
B	943	VAL	-	expression tag	UNP Q8SWT2
B	944	ASP	-	expression tag	UNP Q8SWT2

- Molecule 2 is a RNA chain called RNA (5'-R(P*UP*UP*UP*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	P	0	0	0
			160	72	16	64	8			
2	D	8	Total	C	N	O	P	0	0	0
			160	72	16	64	8			

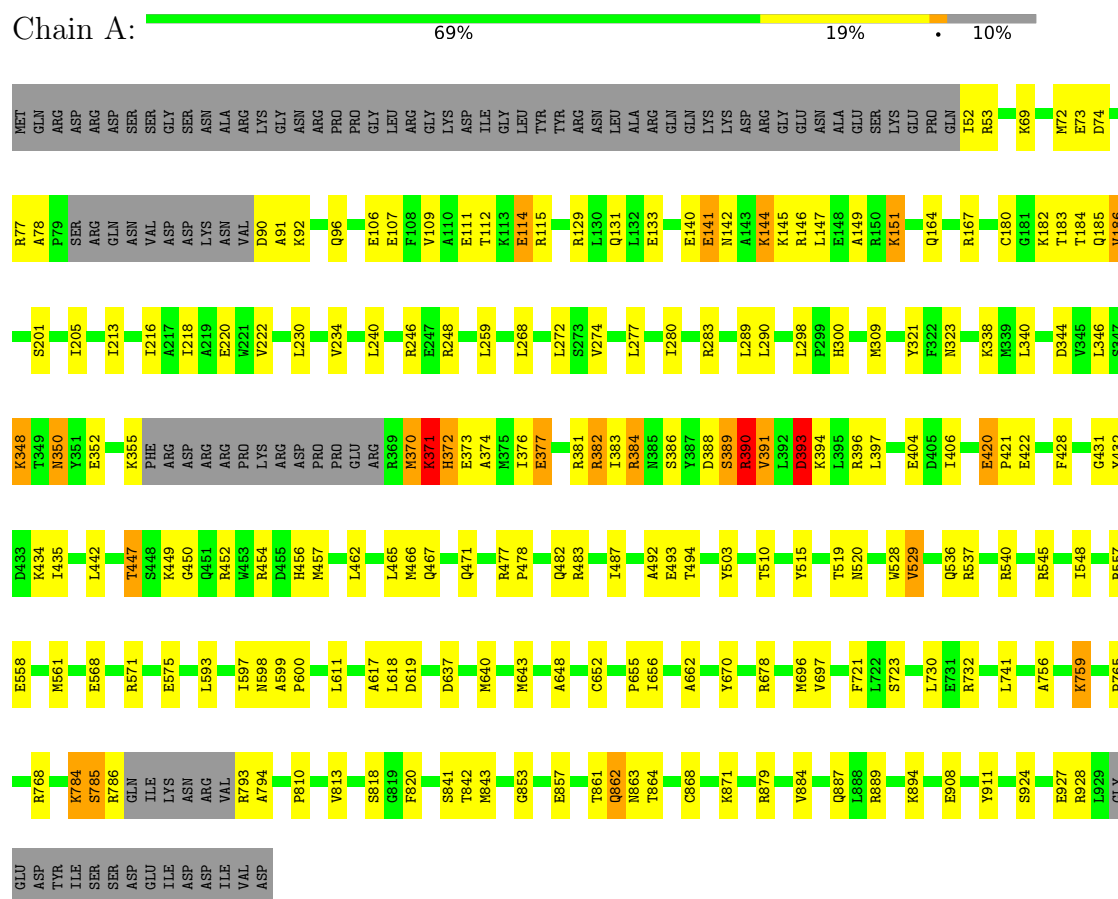
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		
3	B	29	Total	O	0	0
			29	29		
3	C	4	Total	O	0	0
			4	4		
3	D	4	Total	O	0	0
			4	4		

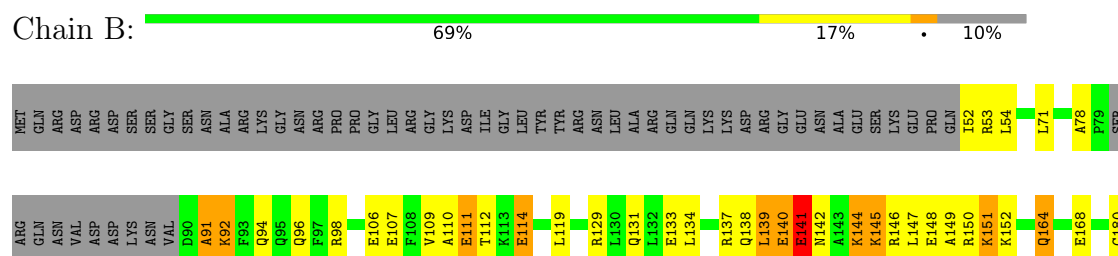
3 Residue-property plots

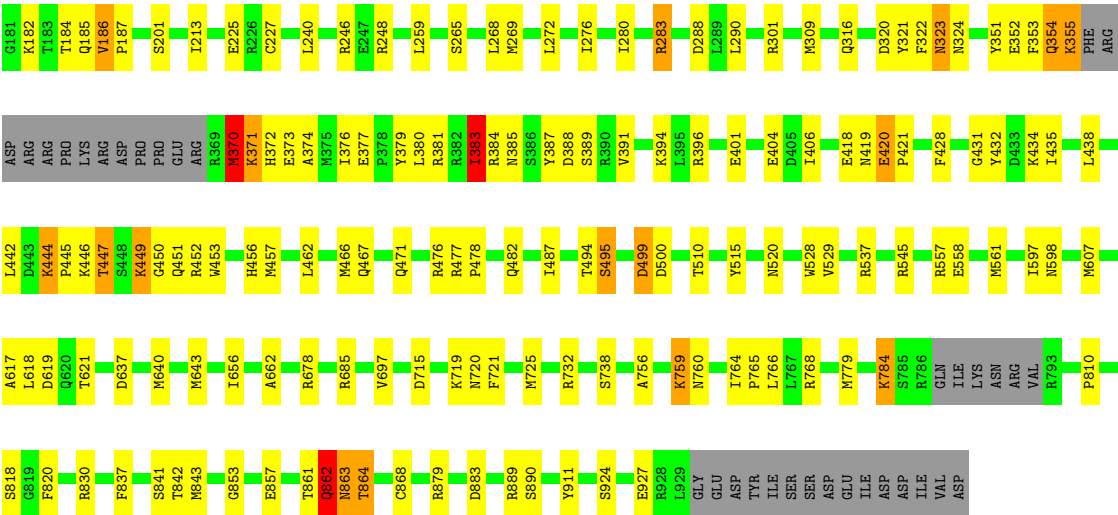
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: CG9323, isoform A



• Molecule 1: CG9323, isoform A

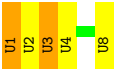
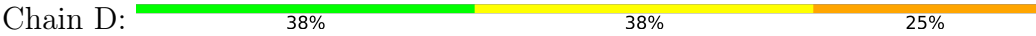




● Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*U)-3')



● Molecule 2: RNA (5'-R(P*UP*UP*UP*UP*UP*UP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	338.33Å 51.53Å 159.90Å 90.00° 118.15° 90.00°	Depositor
Resolution (Å)	47.00 – 2.43 47.00 – 2.43	Depositor EDS
% Data completeness (in resolution range)	90.8 (47.00-2.43) 90.5 (47.00-2.43)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.42Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.212 , 0.278 0.221 , 0.285	Depositor DCC
R_{free} test set	4043 reflections (4.34%)	wwPDB-VP
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.747	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 30.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.469 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14015	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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	0.62	1/6933 (0.0%)	1.01	21/9353 (0.2%)
1	B	0.63	3/6933 (0.0%)	0.99	17/9353 (0.2%)
2	C	0.38	0/175	0.75	0/268
2	D	0.39	0/175	0.74	0/268
All	All	0.62	4/14216 (0.0%)	0.99	38/19242 (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	0	3
1	B	0	4
All	All	0	7

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	864	THR	CA-C	-5.88	1.45	1.53
1	B	383	ILE	CA-CB	5.41	1.60	1.54
1	B	186	VAL	CA-CB	-5.31	1.51	1.54
1	A	186	VAL	CA-CB	-5.04	1.51	1.54

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	141	GLU	N-CA-C	9.62	125.20	107.60
1	A	391	VAL	N-CA-C	-8.66	100.47	112.50
1	A	142	ASN	N-CA-C	-8.42	100.10	110.88
1	B	863	ASN	N-CA-C	-7.06	102.25	111.71
1	B	139	LEU	N-CA-C	-6.80	105.92	114.56
1	A	784	LYS	N-CA-C	6.53	119.33	109.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	145	LYS	N-CA-C	-6.47	103.92	110.97
1	B	91	ALA	N-CA-C	6.38	119.11	108.20
1	A	370	MET	CA-C-N	6.36	133.15	121.70
1	A	370	MET	C-N-CA	6.36	133.15	121.70
1	B	385	ASN	N-CA-C	6.29	118.22	111.36
1	B	862	GLN	CA-CB-CG	6.24	126.58	114.10
1	A	519	THR	N-CA-C	-6.23	105.46	114.12
1	B	111	GLU	N-CA-C	6.01	118.79	111.82
1	B	370	MET	CA-C-N	5.90	132.33	121.70
1	B	370	MET	C-N-CA	5.90	132.33	121.70
1	A	393	ASP	CA-C-N	5.67	128.65	120.38
1	A	393	ASP	C-N-CA	5.67	128.65	120.38
1	B	355	LYS	CA-CB-CG	5.64	125.39	114.10
1	A	91	ALA	N-CA-C	5.61	119.89	111.81
1	A	619	ASP	CA-C-N	-5.57	113.76	122.73
1	A	619	ASP	C-N-CA	-5.57	113.76	122.73
1	B	841	SER	N-CA-C	5.56	116.97	108.52
1	A	841	SER	N-CA-C	5.54	116.94	108.52
1	B	619	ASP	CA-C-N	-5.44	113.94	122.65
1	B	619	ASP	C-N-CA	-5.44	113.94	122.65
1	B	140	GLU	N-CA-C	-5.36	103.42	110.33
1	A	390	ARG	CA-C-N	5.25	129.24	120.64
1	A	390	ARG	C-N-CA	5.25	129.24	120.64
1	A	371	LYS	N-CA-C	5.18	125.49	111.00
1	A	377	GLU	CA-C-N	-5.16	113.72	119.19
1	A	377	GLU	C-N-CA	-5.16	113.72	119.19
1	B	92	LYS	N-CA-C	-5.14	105.85	111.82
1	B	495	SER	N-CA-C	5.14	116.88	111.28
1	B	353	PHE	N-CA-C	5.03	120.31	110.56
1	A	298	LEU	CA-C-N	-5.02	113.76	119.28
1	A	298	LEU	C-N-CA	-5.02	113.76	119.28
1	A	372	HIS	N-CA-C	5.01	118.86	112.34

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	141	GLU	Peptide
1	A	390	ARG	Peptide
1	A	90	ASP	Peptide
1	B	141	GLU	Peptide
1	B	322	PHE	Peptide

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Mol	Chain	Res	Type	Group
1	B	354	GLN	Peptide
1	B	862	GLN	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	6814	0	6913	134	0
1	B	6814	0	6913	130	0
2	C	160	0	81	6	0
2	D	160	0	81	5	0
3	A	30	0	0	3	0
3	B	29	0	0	1	0
3	C	4	0	0	0	0
3	D	4	0	0	1	0
All	All	14015	0	13988	266	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 10.

All (266) 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:D:8:U:O3'	3:D:101:HOH:O	1.84	0.95
1:B:862:GLN:HB3	1:B:864:THR:H	1.33	0.91
1:B:283:ARG:NH2	1:B:598:ASN:OD1	2.02	0.91
1:A:373:GLU:HA	1:A:376:ILE:HG22	1.51	0.90
1:B:323:ASN:OD1	1:B:324:ASN:N	2.04	0.89
1:A:377:GLU:OE1	1:A:396:ARG:NH2	2.08	0.87
1:B:133:GLU:O	1:B:137:ARG:HG3	1.79	0.82
1:B:373:GLU:HA	1:B:376:ILE:HG22	1.59	0.82
1:B:145:LYS:HZ3	1:B:146:ARG:H	1.26	0.81
1:A:390:ARG:CG	1:A:393:ASP:HB3	2.12	0.80
1:A:114:GLU:OE1	1:A:114:GLU:N	2.14	0.79
1:B:477:ARG:HG2	1:B:478:PRO:HD2	1.65	0.79
1:A:393:ASP:CG	1:A:394:LYS:H	1.88	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:GLU:N	1:B:114:GLU:OE1	2.15	0.76
1:A:164:GLN:HE22	1:A:167:ARG:NH1	1.84	0.76
1:B:406:ILE:HD11	1:B:434:LYS:HD3	1.68	0.76
1:B:862:GLN:OE1	1:B:863:ASN:N	2.13	0.75
1:A:389:SER:OG	1:A:390:ARG:N	2.19	0.74
1:B:164:GLN:NE2	1:B:168:GLU:OE2	2.20	0.74
1:B:370:MET:HA	1:B:372:HIS:H	1.52	0.74
1:A:393:ASP:CG	1:A:394:LYS:HG3	2.14	0.73
1:B:144:LYS:HA	1:B:147:LEU:HB3	1.70	0.73
1:A:393:ASP:OD1	1:A:394:LYS:N	2.22	0.72
1:A:390:ARG:HG3	1:A:393:ASP:HB3	1.69	0.72
1:A:52:ILE:HD11	1:A:861:THR:HG21	1.71	0.71
1:B:765:PRO:HG2	1:B:927:GLU:HG2	1.73	0.71
1:B:418:GLU:HG2	1:B:453:TRP:HZ2	1.54	0.71
1:A:283:ARG:O	3:A:1001:HOH:O	2.08	0.70
1:A:449:LYS:HA	1:A:452:ARG:HD3	1.73	0.69
1:B:924:SER:O	1:B:927:GLU:HG3	1.92	0.69
1:A:810:PRO:HB2	2:C:1:U:C6	2.26	0.69
1:A:447:THR:HG22	1:A:450:GLY:H	1.58	0.69
1:B:853:GLY:O	1:B:889:ARG:NH2	2.26	0.68
1:B:379:TYR:O	1:B:383:ILE:HG22	1.94	0.68
1:B:381:ARG:O	1:B:384:ARG:HG3	1.94	0.68
1:A:107:GLU:O	1:A:111:GLU:HG3	1.94	0.67
1:B:146:ARG:NH1	1:B:227:CYS:SG	2.68	0.67
1:B:351:TYR:O	1:B:394:LYS:NZ	2.24	0.67
1:A:131:GLN:OE1	3:A:1002:HOH:O	2.11	0.67
1:B:449:LYS:HB2	1:B:452:ARG:NH1	2.10	0.66
1:A:146:ARG:HA	1:A:149:ALA:HB3	1.78	0.66
1:A:862:GLN:OE1	1:A:863:ASN:N	2.28	0.66
1:B:370:MET:HA	1:B:372:HIS:N	2.09	0.66
1:B:146:ARG:HD3	1:B:225:GLU:HA	1.78	0.65
1:A:477:ARG:HG3	1:A:477:ARG:HH11	1.62	0.65
1:B:240:LEU:HD13	1:B:732:ARG:HB3	1.78	0.64
1:B:129:ARG:O	1:B:133:GLU:HG3	1.97	0.64
1:A:637:ASP:HB2	1:A:640:MET:HE2	1.80	0.64
1:A:924:SER:O	1:A:927:GLU:HG3	1.98	0.63
1:B:449:LYS:HB2	1:B:452:ARG:HH12	1.61	0.63
1:B:857:GLU:HG2	1:B:868:CYS:SG	2.38	0.63
1:B:418:GLU:HG2	1:B:453:TRP:CZ2	2.34	0.63
1:A:259:LEU:HD23	1:A:290:LEU:HD21	1.82	0.62
1:B:447:THR:HG22	1:B:450:GLY:H	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:810:PRO:HB2	2:C:1:U:H6	1.63	0.62
1:A:129:ARG:O	1:A:133:GLU:HG3	2.00	0.61
1:B:377:GLU:OE1	1:B:396:ARG:NH1	2.30	0.61
1:B:467:GLN:OE1	1:B:471:GLN:NE2	2.27	0.61
1:A:515:TYR:OH	1:A:520:ASN:OD1	2.16	0.60
1:A:164:GLN:HE22	1:A:167:ARG:HH12	1.49	0.60
1:A:447:THR:HG22	1:A:450:GLY:N	2.15	0.60
1:B:499:ASP:OD1	1:B:499:ASP:N	2.32	0.60
1:A:406:ILE:HD11	1:A:434:LYS:HD3	1.83	0.60
1:B:112:THR:HG21	1:B:268:LEU:HD13	1.84	0.60
1:B:147:LEU:O	1:B:151:LYS:HG2	2.01	0.59
1:B:685:ARG:NH2	1:B:720:ASN:OD1	2.36	0.59
1:A:370:MET:HA	1:A:372:HIS:N	2.18	0.58
1:B:107:GLU:O	1:B:111:GLU:HG3	2.02	0.58
1:B:146:ARG:HA	1:B:149:ALA:HB3	1.85	0.58
1:A:240:LEU:HD13	1:A:732:ARG:HB3	1.86	0.57
1:A:431:GLY:O	1:A:435:ILE:HG13	2.04	0.57
1:B:141:GLU:OE1	1:B:144:LYS:HG2	2.05	0.57
1:A:529:VAL:O	3:A:1003:HOH:O	2.17	0.56
1:A:756:ALA:O	1:A:759:LYS:HE3	2.06	0.56
1:A:283:ARG:NH2	1:A:598:ASN:OD1	2.33	0.56
1:A:467:GLN:OE1	1:A:471:GLN:NE2	2.35	0.56
1:B:52:ILE:C	1:B:53:ARG:HD3	2.31	0.56
1:B:91:ALA:HB1	1:B:94:GLN:H	1.71	0.56
1:B:148:GLU:O	1:B:152:LYS:HG3	2.05	0.56
1:A:449:LYS:HA	1:A:452:ARG:CD	2.36	0.55
1:A:759:LYS:NZ	1:A:908:GLU:OE1	2.39	0.55
1:B:431:GLY:O	1:B:435:ILE:HG13	2.07	0.55
1:B:383:ILE:HD11	1:B:387:TYR:CD2	2.41	0.55
1:A:69:LYS:O	1:A:73:GLU:HG3	2.07	0.55
1:A:381:ARG:O	1:A:383:ILE:N	2.40	0.54
1:A:370:MET:SD	1:A:373:GLU:HG2	2.48	0.54
1:B:756:ALA:O	1:B:759:LYS:NZ	2.31	0.54
1:B:142:ASN:HB2	1:B:145:LYS:HD3	1.90	0.54
1:B:377:GLU:O	1:B:381:ARG:HG3	2.08	0.54
1:B:445:PRO:HG2	1:B:451:GLN:HG2	1.89	0.54
1:B:447:THR:HG22	1:B:449:LYS:H	1.73	0.54
1:A:390:ARG:HG2	1:A:393:ASP:HB3	1.86	0.54
1:B:78:ALA:HB2	1:B:911:TYR:CG	2.44	0.54
1:A:109:VAL:HG22	1:A:268:LEU:HD21	1.90	0.53
1:A:390:ARG:HG3	1:A:393:ASP:CB	2.36	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:LYS:O	1:B:96:GLN:HG3	2.08	0.53
1:B:820:PHE:HB3	1:B:843:MET:SD	2.48	0.53
1:A:388:ASP:OD1	1:A:389:SER:N	2.42	0.53
1:A:492:ALA:O	1:A:540:ARG:NH1	2.42	0.53
1:B:477:ARG:CG	1:B:478:PRO:HD2	2.36	0.52
1:A:765:PRO:HG2	1:A:927:GLU:HG2	1.90	0.52
1:B:201:SER:O	1:B:248:ARG:NH1	2.41	0.52
1:A:340:LEU:HG	1:A:548:ILE:HD11	1.90	0.52
1:A:390:ARG:HE	1:A:393:ASP:HB3	1.73	0.52
1:A:393:ASP:OD2	1:A:394:LYS:HE3	2.09	0.52
1:A:112:THR:HG21	1:A:268:LEU:HD13	1.91	0.52
1:B:637:ASP:HB2	1:B:640:MET:HE2	1.90	0.52
1:A:853:GLY:O	1:A:889:ARG:NH2	2.43	0.52
1:B:109:VAL:HG22	1:B:268:LEU:HD21	1.93	0.51
1:B:810:PRO:HG2	2:D:1:U:H1'	1.93	0.51
1:A:92:LYS:O	1:A:96:GLN:HG3	2.11	0.51
1:A:140:GLU:HG2	1:A:141:GLU:N	2.24	0.51
1:A:768:ARG:HD3	1:A:843:MET:O	2.11	0.50
1:B:134:LEU:O	1:B:138:GLN:HG3	2.11	0.50
1:A:820:PHE:HB3	1:A:843:MET:SD	2.52	0.49
1:A:52:ILE:C	1:A:53:ARG:HD3	2.37	0.49
1:A:467:GLN:HG3	1:A:670:TYR:CZ	2.48	0.49
1:B:371:LYS:O	1:B:374:ALA:N	2.45	0.49
1:B:131:GLN:OE1	3:B:1001:HOH:O	2.20	0.49
1:A:78:ALA:HB2	1:A:911:TYR:CD2	2.47	0.49
1:A:393:ASP:OD1	1:A:394:LYS:HG3	2.13	0.49
1:A:74:ASP:O	1:A:77:ARG:HG2	2.12	0.49
1:A:477:ARG:HG3	1:A:477:ARG:NH1	2.25	0.48
1:B:456:HIS:HA	1:B:482:GLN:HG2	1.95	0.48
1:B:617:ALA:O	1:B:618:LEU:HD23	2.14	0.48
1:B:810:PRO:HB2	2:D:1:U:C6	2.48	0.48
1:A:53:ARG:HD3	1:A:53:ARG:N	2.29	0.48
1:A:180:CYS:SG	1:A:182:LYS:HG2	2.53	0.48
1:A:371:LYS:O	1:A:374:ALA:N	2.46	0.48
1:A:421:PRO:HB2	1:A:545:ARG:HH22	1.78	0.47
1:B:442:LEU:HD22	1:B:457:MET:HE1	1.96	0.47
1:A:186:VAL:HG21	1:A:309:MET:CE	2.44	0.47
1:B:145:LYS:HZ3	1:B:146:ARG:HG3	1.79	0.47
2:D:3:U:H2'	2:D:4:U:O4'	2.15	0.47
1:B:764:ILE:O	1:B:768:ARG:HG3	2.15	0.47
1:A:785:SER:OG	1:A:786:ARG:N	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:432:TYR:CE2	1:B:466:MET:HE2	2.50	0.47
1:B:186:VAL:HG21	1:B:309:MET:CE	2.45	0.47
1:A:558:GLU:HA	1:A:561:MET:SD	2.55	0.47
1:B:184:THR:OG1	1:B:185:GLN:NE2	2.39	0.47
1:A:390:ARG:HE	1:A:393:ASP:CB	2.28	0.47
1:B:419:ASN:O	1:B:420:GLU:HB2	2.15	0.47
1:B:180:CYS:SG	1:B:182:LYS:HG2	2.54	0.46
1:B:428:PHE:CE2	1:B:537:ARG:HG2	2.50	0.46
1:B:52:ILE:HD11	1:B:861:THR:HG21	1.96	0.46
1:B:186:VAL:HG21	1:B:309:MET:HE1	1.98	0.46
1:B:421:PRO:HB2	1:B:545:ARG:HH22	1.79	0.46
1:B:447:THR:CG2	1:B:449:LYS:H	2.28	0.46
1:A:432:TYR:CE2	1:A:466:MET:HE2	2.51	0.46
1:B:381:ARG:O	1:B:383:ILE:N	2.49	0.46
1:B:528:TRP:CZ2	1:B:557:ARG:HD3	2.49	0.46
1:B:146:ARG:O	1:B:150:ARG:N	2.38	0.46
1:A:377:GLU:O	1:A:381:ARG:HG3	2.15	0.46
1:B:283:ARG:NH1	1:B:288:ASP:OD2	2.49	0.46
1:B:438:LEU:HD23	1:B:487:ILE:HG23	1.98	0.45
1:A:390:ARG:HG3	1:A:393:ASP:CA	2.46	0.45
1:B:371:LYS:NZ	1:B:557:ARG:HH22	2.14	0.45
1:A:394:LYS:HA	1:A:397:LEU:HG	1.98	0.45
1:A:617:ALA:O	1:A:618:LEU:HD23	2.17	0.45
1:B:376:ILE:O	1:B:380:LEU:HG	2.15	0.45
1:A:810:PRO:HB2	2:C:1:U:C5	2.51	0.45
1:B:383:ILE:HG12	1:B:383:ILE:O	2.16	0.45
1:B:494:THR:OG1	1:B:495:SER:N	2.50	0.45
1:A:205:ILE:HG12	1:A:274:VAL:HB	1.97	0.45
1:A:528:TRP:CE2	1:A:557:ARG:HD3	2.52	0.45
1:B:142:ASN:CB	1:B:145:LYS:HD3	2.46	0.45
1:B:447:THR:HG22	1:B:449:LYS:N	2.31	0.45
1:B:879:ARG:NH1	1:B:883:ASP:OD1	2.42	0.45
1:A:793:ARG:CG	1:A:794:ALA:N	2.80	0.45
1:A:857:GLU:HB2	1:A:871:LYS:HE2	1.99	0.45
1:A:201:SER:O	1:A:248:ARG:NH1	2.45	0.45
1:A:643:MET:HE3	1:A:656:ILE:HG23	1.98	0.45
1:A:662:ALA:HB2	1:A:697:VAL:HG11	1.99	0.45
1:A:183:THR:HG21	1:A:218:ILE:HD13	1.99	0.44
1:A:477:ARG:HB3	1:A:478:PRO:HD2	1.99	0.44
1:A:670:TYR:CZ	1:A:723:SER:HB2	2.52	0.44
1:B:140:GLU:HG2	1:B:141:GLU:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:528:TRP:CZ2	1:A:557:ARG:HD3	2.53	0.44
1:B:71:LEU:HA	1:B:71:LEU:HD23	1.72	0.44
1:B:111:GLU:HA	1:B:114:GLU:CD	2.43	0.44
1:B:447:THR:HG22	1:B:450:GLY:N	2.29	0.44
1:A:696:MET:HE2	1:A:813:VAL:CG1	2.47	0.44
1:B:138:GLN:O	1:B:139:LEU:HD23	2.17	0.44
1:B:388:ASP:OD1	1:B:389:SER:N	2.51	0.44
1:B:421:PRO:HB2	1:B:545:ARG:NH2	2.32	0.44
1:B:784:LYS:HA	1:B:784:LYS:HD2	1.81	0.44
1:B:515:TYR:OH	1:B:520:ASN:OD1	2.24	0.44
1:B:265:SER:HG	2:D:8:U:H3	1.62	0.44
1:B:316:GLN:NE2	1:B:320:ASP:OD1	2.42	0.44
1:A:447:THR:CG2	1:A:449:LYS:H	2.30	0.43
1:B:144:LYS:HA	1:B:147:LEU:CB	2.45	0.43
1:B:558:GLU:HA	1:B:561:MET:SD	2.58	0.43
1:A:346:LEU:HD23	1:A:346:LEU:HA	1.82	0.43
1:A:370:MET:HA	1:A:372:HIS:H	1.83	0.43
1:A:421:PRO:HB2	1:A:545:ARG:NH2	2.33	0.43
1:A:652:CYS:O	1:A:655:PRO:HD2	2.18	0.43
1:A:810:PRO:CB	2:C:1:U:H6	2.30	0.43
1:B:768:ARG:HD3	1:B:843:MET:O	2.17	0.43
1:A:147:LEU:O	1:A:151:LYS:HB2	2.18	0.43
1:A:428:PHE:CE2	1:A:537:ARG:HG2	2.53	0.43
1:B:269:MET:O	1:B:301:ARG:NH1	2.50	0.43
1:A:184:THR:OG1	1:A:185:GLN:NE2	2.46	0.43
1:A:420:GLU:HA	1:A:421:PRO:HD3	1.75	0.43
1:B:187:PRO:HG3	1:B:276:ILE:HD12	2.00	0.43
1:B:401:GLU:OE1	1:B:557:ARG:HD2	2.19	0.43
1:A:861:THR:HG23	1:A:862:GLN:N	2.34	0.43
1:A:862:GLN:CD	1:A:863:ASN:N	2.77	0.43
1:B:52:ILE:HD11	1:B:861:THR:OG1	2.19	0.43
1:A:72:MET:HE2	1:A:884:VAL:HG13	2.00	0.43
1:A:216:ILE:O	1:A:220:GLU:HG3	2.19	0.43
1:B:830:ARG:HA	1:B:837:PHE:O	2.19	0.43
1:B:662:ALA:HB2	1:B:697:VAL:HG11	2.01	0.43
1:B:678:ARG:HG2	1:B:721:PHE:HE2	1.83	0.43
1:A:862:GLN:OE1	1:A:864:THR:HG23	2.19	0.42
1:A:78:ALA:HB2	1:A:911:TYR:CG	2.54	0.42
1:B:760:ASN:HB3	1:B:766:LEU:HD23	2.02	0.42
1:A:381:ARG:C	1:A:383:ILE:N	2.78	0.42
1:B:145:LYS:HE2	1:B:146:ARG:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:GLU:HG2	1:A:536:GLN:HG2	2.01	0.42
1:B:607:MET:HE3	1:B:607:MET:HB2	1.80	0.42
1:B:779:MET:HE3	1:B:779:MET:HB2	1.93	0.42
1:B:107:GLU:O	1:B:110:ALA:HB3	2.20	0.42
1:B:377:GLU:O	1:B:377:GLU:HG3	2.19	0.42
1:A:289:LEU:HD23	1:A:289:LEU:HA	1.92	0.42
1:A:421:PRO:HG2	1:A:545:ARG:HH22	1.84	0.42
1:A:277:LEU:HD21	1:A:290:LEU:HD23	2.02	0.42
1:A:503:TYR:HD1	1:A:548:ILE:HG22	1.84	0.42
1:A:442:LEU:HB3	1:A:457:MET:HE1	2.01	0.42
1:B:246:ARG:HG3	1:B:248:ARG:O	2.20	0.42
1:A:422:GLU:HA	1:A:483:ARG:HH12	1.85	0.41
1:A:857:GLU:HB3	1:A:868:CYS:SG	2.60	0.41
1:A:381:ARG:O	1:A:384:ARG:HG3	2.20	0.41
1:A:599:ALA:HA	1:A:600:PRO:HD3	1.91	0.41
1:B:500:ASP:OD1	1:B:500:ASP:N	2.54	0.41
1:A:246:ARG:HG3	1:A:248:ARG:O	2.20	0.41
1:A:456:HIS:C	1:A:482:GLN:HG2	2.46	0.41
1:A:862:GLN:O	1:A:863:ASN:ND2	2.54	0.41
1:B:54:LEU:HA	1:B:54:LEU:HD23	1.88	0.41
1:B:110:ALA:O	1:B:114:GLU:OE1	2.39	0.41
1:B:715:ASP:OD2	1:B:719:LYS:NZ	2.32	0.41
1:A:730:LEU:HD23	1:A:730:LEU:HA	1.80	0.41
1:B:145:LYS:NZ	1:B:146:ARG:H	2.08	0.41
1:A:321:TYR:CD1	1:A:597:ILE:HG12	2.56	0.41
1:B:321:TYR:CD1	1:B:597:ILE:HG12	2.55	0.41
1:B:421:PRO:HG2	1:B:545:ARG:HH22	1.84	0.41
1:A:394:LYS:H	1:A:394:LYS:HG3	1.71	0.41
1:A:503:TYR:CD1	1:A:548:ILE:HG22	2.56	0.41
1:A:678:ARG:HG2	1:A:721:PHE:HE2	1.85	0.41
1:A:344:ASP:O	1:A:348:LYS:HG3	2.21	0.41
1:A:648:ALA:HB2	1:A:741:LEU:HD23	2.03	0.41
1:A:810:PRO:HG2	2:C:1:U:H1'	2.03	0.41
1:B:462:LEU:HD12	1:B:462:LEU:HA	1.91	0.41
1:A:230:LEU:HD23	1:A:230:LEU:HA	1.80	0.41
1:B:444:LYS:O	1:B:444:LYS:HG2	2.21	0.41
1:A:222:VAL:HG12	1:A:234:VAL:HG21	2.02	0.40
1:A:465:LEU:HD23	2:C:2:U:O2	2.20	0.40
1:B:643:MET:HE3	1:B:656:ILE:HG23	2.04	0.40
1:B:145:LYS:NZ	1:B:146:ARG:HG3	2.36	0.40
1:B:259:LEU:HD23	1:B:290:LEU:HD21	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:ARG:O	1:A:382:ARG:C	2.65	0.40
1:A:462:LEU:N	1:A:487:ILE:O	2.54	0.40
1:A:568:GLU:HA	1:A:571:ARG:HE	1.87	0.40
1:A:575:GLU:HG2	1:A:611:LEU:HD22	2.04	0.40
1:B:447:THR:O	1:B:451:GLN:HG3	2.21	0.40
1:A:593:LEU:HA	1:A:593:LEU:HD23	1.80	0.40
1:B:372:HIS:O	1:B:372:HIS:CG	2.74	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	841/944 (89%)	792 (94%)	38 (4%)	11 (1%)	9	13
1	B	841/944 (89%)	793 (94%)	41 (5%)	7 (1%)	16	23
All	All	1682/1888 (89%)	1585 (94%)	79 (5%)	18 (1%)	11	16

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	350	ASN
1	A	371	LYS
1	A	389	SER
1	A	785	SER
1	B	323	ASN
1	B	371	LYS
1	B	862	GLN
1	A	144	LYS
1	A	323	ASN
1	A	393	ASP
1	A	382	ARG

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Mol	Chain	Res	Type
1	A	420	GLU
1	B	354	GLN
1	B	144	LYS
1	B	420	GLU
1	A	529	VAL
1	A	862	GLN
1	B	529	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	757/842 (90%)	727 (96%)	30 (4%)	28	45
1	B	757/842 (90%)	725 (96%)	32 (4%)	26	43
All	All	1514/1684 (90%)	1452 (96%)	62 (4%)	27	44

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	GLU
1	A	114	GLU
1	A	115	ARG
1	A	144	LYS
1	A	151	LYS
1	A	213	ILE
1	A	272	LEU
1	A	280	ILE
1	A	300	HIS
1	A	338	LYS
1	A	348	LYS
1	A	350	ASN
1	A	352	GLU
1	A	355	LYS
1	A	384	ARG
1	A	386	SER

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Mol	Chain	Res	Type
1	A	391	VAL
1	A	404	GLU
1	A	447	THR
1	A	454	ARG
1	A	494	THR
1	A	510	THR
1	A	759	LYS
1	A	784	LYS
1	A	818	SER
1	A	842	THR
1	A	879	ARG
1	A	887	GLN
1	A	894	LYS
1	A	928	ARG
1	B	98	ARG
1	B	106	GLU
1	B	114	GLU
1	B	119	LEU
1	B	145	LYS
1	B	151	LYS
1	B	164	GLN
1	B	213	ILE
1	B	272	LEU
1	B	280	ILE
1	B	283	ARG
1	B	352	GLU
1	B	355	LYS
1	B	370	MET
1	B	383	ILE
1	B	391	VAL
1	B	404	GLU
1	B	444	LYS
1	B	446	LYS
1	B	447	THR
1	B	449	LYS
1	B	476	ARG
1	B	499	ASP
1	B	510	THR
1	B	621	THR
1	B	725	MET
1	B	738	SER
1	B	759	LYS

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Mol	Chain	Res	Type
1	B	784	LYS
1	B	818	SER
1	B	842	THR
1	B	890	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	GLN
1	A	237	GLN
1	A	264	GLN
1	A	300	HIS
1	A	863	ASN
1	A	909	ASN
1	B	232	ASN
1	B	261	GLN
1	B	300	HIS
1	B	419	ASN
1	B	456	HIS
1	B	728	GLN
1	B	807	ASN
1	B	914	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	8/8 (100%)	2 (25%)	1 (12%)
2	D	8/8 (100%)	2 (25%)	1 (12%)
All	All	16/16 (100%)	4 (25%)	2 (12%)

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	2	U
2	C	3	U
2	D	2	U
2	D	3	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	C	1	U
2	D	1	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](#)

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	849/944 (89%)	-1.10	0 100 100	37, 57, 90, 112	0
1	B	849/944 (89%)	-1.11	0 100 100	36, 57, 90, 108	0
2	C	8/8 (100%)	-1.64	0 100 100	43, 56, 71, 83	0
2	D	8/8 (100%)	-1.60	0 100 100	44, 58, 73, 83	0
All	All	1714/1904 (90%)	-1.11	0 100 100	36, 57, 90, 112	0

There are no RSRZ outliers to report.

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.