



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:22 PM UTC

PDB ID : 7XX5 / pdb_00007xx5
Title : Crystal Structure of Nucleosome-H1.3 Linker Histone Assembly (sticky-169a DNA fragment)
Authors : Adhireksan, Z.; Qiuye, B.; Lee, P.L.; Sharma, D.; Padavattan, S.; Davey, C.A.
Deposited on : 2022-05-28
Resolution : 3.19 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

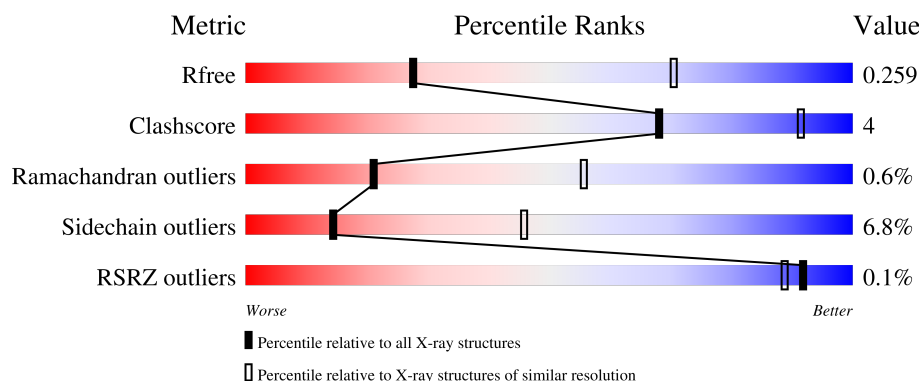
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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.



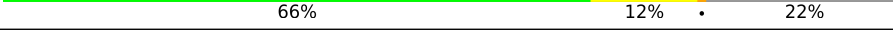
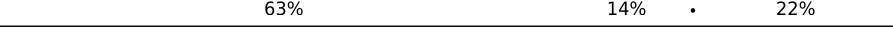
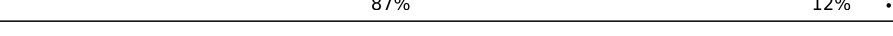
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	138	
1	E	138	
1	K	138	
1	O	138	
2	B	105	

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Mol	Chain	Length	Quality of chain
2	F	105	
2	L	105	
2	P	105	
3	C	132	
3	G	132	
3	M	132	
3	Q	132	
4	D	128	
4	H	128	
4	N	128	
4	R	128	
5	I	169	
5	S	169	
6	J	169	
6	T	169	
7	U	222	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 26617 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 Histone H3.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			816	514	158	140	4			
1	E	99	Total	C	N	O	S	0	0	0
			816	514	158	140	4			
1	K	99	Total	C	N	O	S	0	0	0
			816	514	158	140	4			
1	O	100	Total	C	N	O	S	0	0	0
			825	520	160	141	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P68431
A	-1	SER	-	expression tag	UNP P68431
E	-2	GLY	-	expression tag	UNP P68431
E	-1	SER	-	expression tag	UNP P68431
K	-2	GLY	-	expression tag	UNP P68431
K	-1	SER	-	expression tag	UNP P68431
O	-2	GLY	-	expression tag	UNP P68431
O	-1	SER	-	expression tag	UNP P68431

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	80	Total	C	N	O	S	0	0	0
			638	401	125	111	1			
2	F	80	Total	C	N	O	S	0	0	0
			638	401	125	111	1			
2	L	80	Total	C	N	O	S	0	0	0
			638	401	125	111	1			
2	P	80	Total	C	N	O	S	0	0	0
			638	401	125	111	1			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP P62805
B	-1	SER	-	expression tag	UNP P62805
F	-2	GLY	-	expression tag	UNP P62805
F	-1	SER	-	expression tag	UNP P62805
L	-2	GLY	-	expression tag	UNP P62805
L	-1	SER	-	expression tag	UNP P62805
P	-2	GLY	-	expression tag	UNP P62805
P	-1	SER	-	expression tag	UNP P62805

- Molecule 3 is a protein called Histone H2A type 1-B/E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	105	Total	C	N	O	0	0	0
			810	511	158	141			
3	G	103	Total	C	N	O	0	0	0
			796	502	155	139			
3	M	105	Total	C	N	O	0	0	0
			810	511	158	141			
3	Q	104	Total	C	N	O	0	0	0
			805	508	157	140			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP P04908
C	-1	SER	-	expression tag	UNP P04908
G	-2	GLY	-	expression tag	UNP P04908
G	-1	SER	-	expression tag	UNP P04908
M	-2	GLY	-	expression tag	UNP P04908
M	-1	SER	-	expression tag	UNP P04908
Q	-2	GLY	-	expression tag	UNP P04908
Q	-1	SER	-	expression tag	UNP P04908

- Molecule 4 is a protein called Histone H2B type 1-J.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	100	Total	C	N	O S	0	0	0
			788	494	147	145 2			
4	H	100	Total	C	N	O S	0	0	0
			788	494	147	145 2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	N	100	Total	C	N	O	S	0	0	0
			788	494	147	145	2			
4	R	100	Total	C	N	O	S	0	0	0
			788	494	147	145	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	GLY	-	expression tag	UNP P06899
D	-1	SER	-	expression tag	UNP P06899
H	-2	GLY	-	expression tag	UNP P06899
H	-1	SER	-	expression tag	UNP P06899
N	-2	GLY	-	expression tag	UNP P06899
N	-1	SER	-	expression tag	UNP P06899
R	-2	GLY	-	expression tag	UNP P06899
R	-1	SER	-	expression tag	UNP P06899

- Molecule 5 is a DNA chain called DNA (169-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	169	Total	C	N	O	P	0	0	0
			3462	1646	637	1011	168			
5	S	169	Total	C	N	O	P	0	0	0
			3462	1646	637	1011	168			

- Molecule 6 is a DNA chain called DNA (169-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	J	169	Total	C	N	O	P	0	0	0
			3461	1646	634	1013	168			
6	T	169	Total	C	N	O	P	0	0	0
			3461	1646	634	1013	168			

- Molecule 7 is a protein called Histone H1.3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	U	78	Total	C	N	O	0	0	0
			560	351	104	105			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
U	0	SER	-	expression tag	UNP P16402

- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

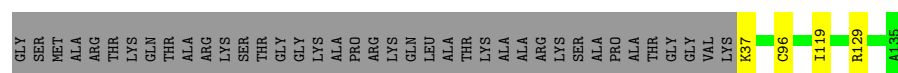
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	G	1	Total Ca 1 1	0	0
8	I	3	Total Ca 3 3	0	0
8	J	1	Total Ca 1 1	0	0
8	Q	3	Total Ca 3 3	0	0
8	S	2	Total Ca 2 2	0	0
8	T	3	Total Ca 3 3	0	0

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.

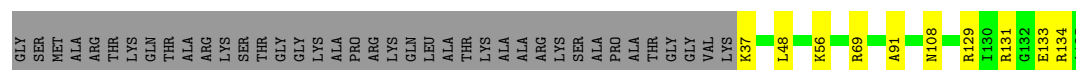
- Molecule 1: Histone H3.1

Chain A: 



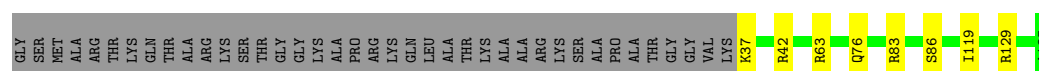
- Molecule 1: Histone H3.1

Chain E: 



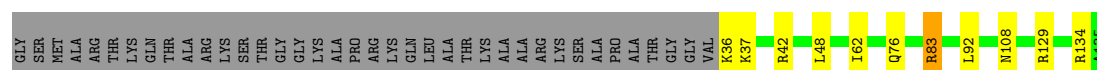
- Molecule 1: Histone H3.1

Chain K: 



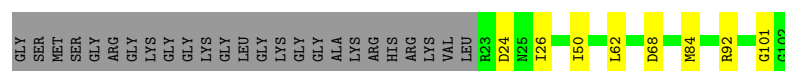
- Molecule 1: Histone H3.1

Chain O: 



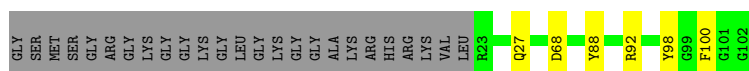
- Molecule 2: Histone H4

Chain B: 

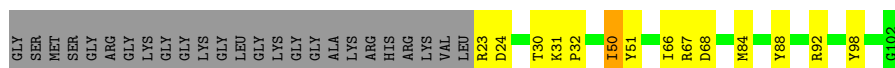


- Molecule 2: Histone H4

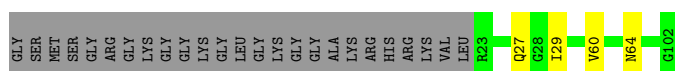
Chain F: 



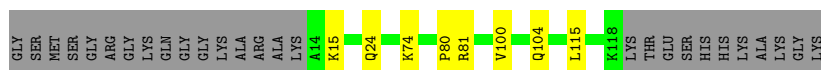
• Molecule 2: Histone H4



• Molecule 2: Histone H4



• Molecule 3: Histone H2A type 1-B/E



• Molecule 3: Histone H2A type 1-B/E



• Molecule 3: Histone H2A type 1-B/E



• Molecule 3: Histone H2A type 1-B/E



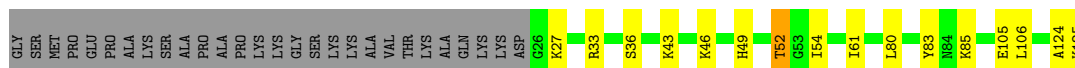
• Molecule 4: Histone H2B type 1-J





- Molecule 4: Histone H2B type 1-J

Chain H: 66% 12% 22%



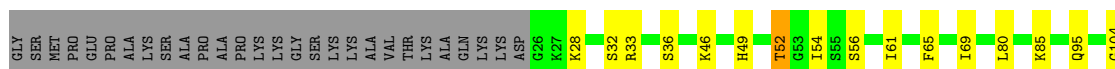
- Molecule 4: Histone H2B type 1-J

Chain N: 67% 11% 22%



- Molecule 4: Histone H2B type 1-J

Chain R: 63% 14% 22%



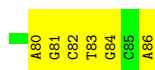
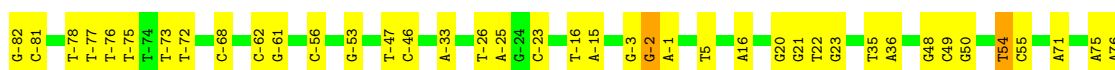
- Molecule 5: DNA (169-MER)

Chain I: 84% 14% 2%

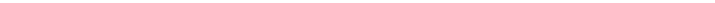


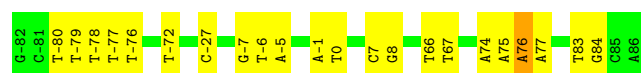
- Molecule 5: DNA (169-MER)

Chain S: 73% 26% 1%

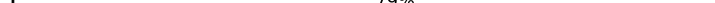


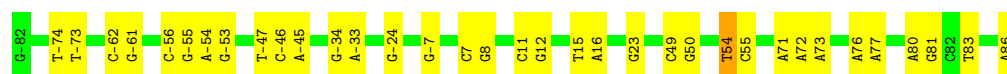
- Molecule 6: DNA (169-MER)

Chain J:  87% 12%

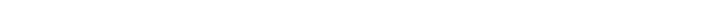


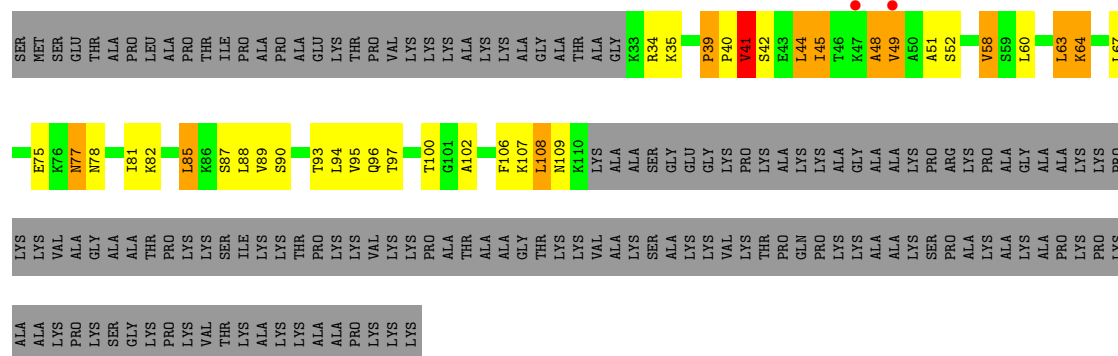
- Molecule 6: DNA (169-MER)

Chain T:  79% 20%



- Molecule 7: Histone H1.3

Chain U:  %



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	105.07Å 101.74Å 218.40Å 90.00° 103.52° 90.00°	Depositor
Resolution (Å)	46.57 – 3.19 46.57 – 3.19	Depositor EDS
% Data completeness (in resolution range)	98.4 (46.57-3.19) 98.4 (46.57-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.203 , 0.259 0.206 , 0.259	Depositor DCC
R_{free} test set	3677 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	102.2	Xtriage
Anisotropy	0.291	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 73.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.059 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	26617	wwPDB-VP
Average B, all atoms (Å ²)	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0/828	0.91	0/1109
1	E	0.63	0/828	0.96	0/1109
1	K	0.73	0/828	1.01	0/1109
1	O	0.77	0/837	0.99	0/1120
2	B	0.61	0/645	0.90	0/862
2	F	0.62	0/645	0.94	0/862
2	L	0.70	0/645	0.91	0/862
2	P	0.82	0/645	0.98	0/862
3	C	0.63	0/820	0.94	0/1107
3	G	0.58	0/806	0.92	0/1089
3	M	0.72	0/820	1.01	0/1107
3	Q	0.73	0/815	0.97	1/1100 (0.1%)
4	D	0.66	0/799	0.97	0/1067
4	H	0.61	0/799	0.94	0/1067
4	N	0.72	0/799	0.98	0/1067
4	R	0.74	0/799	1.00	0/1067
5	I	0.35	0/3884	0.79	4/5993 (0.1%)
5	S	0.43	1/3884 (0.0%)	0.86	4/5993 (0.1%)
6	J	0.34	0/3882	0.80	4/5990 (0.1%)
6	T	0.41	0/3882	0.84	4/5990 (0.1%)
7	U	1.26	2/563 (0.4%)	1.41	6/750 (0.8%)
All	All	0.56	3/28453 (0.0%)	0.90	23/41282 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	U	41	VAL	CA-C	5.53	1.59	1.52
7	U	42	SER	N-CA	5.42	1.53	1.46
5	S	5	DT	O3'-P	-5.02	1.53	1.61

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	80	DA	C2'-C3'-O3'	8.70	124.54	111.50
7	U	41	VAL	N-CA-C	7.41	124.74	109.34
5	I	56	DG	C2'-C3'-O3'	-7.15	100.78	111.50
6	J	77	DA	C2'-C3'-O3'	7.13	122.19	111.50
7	U	85	LEU	N-CA-C	6.97	119.52	111.02
6	J	76	DA	C2'-C3'-O3'	6.70	121.54	111.50
7	U	35	LYS	N-CA-C	6.58	119.73	111.24
6	T	54	DT	C4'-C3'-O3'	6.40	119.61	110.00
6	J	-72	DT	C2'-C3'-O3'	6.35	121.03	111.50
7	U	39	PRO	N-CA-C	6.12	118.17	110.70
6	T	-56	DC	C2'-C3'-O3'	-6.00	102.51	111.50
5	I	79	DA	C2'-C3'-O3'	5.96	120.44	111.50
6	T	-55	DG	C2'-C3'-O3'	-5.80	102.80	111.50
5	S	54	DT	C4'-C3'-O3'	5.76	118.64	110.00
5	S	-2	DG	C2'-C3'-O3'	-5.70	102.94	111.50
5	I	81	DG	C2'-C3'-O3'	5.65	119.97	111.50
5	S	16	DA	C2'-C3'-O3'	-5.53	103.20	111.50
5	S	5	DT	C2'-C3'-O3'	-5.47	103.29	111.50
6	J	-27	DC	C2'-C3'-O3'	-5.36	103.46	111.50
6	T	23	DG	C2'-C3'-O3'	5.31	119.47	111.50
7	U	58	VAL	N-CA-C	5.21	117.19	109.17
7	U	48	ALA	N-CA-C	5.20	116.75	111.14
3	Q	49	VAL	N-CA-C	-5.18	105.55	110.42

There are no chirality outliers.

There are no planarity outliers.

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	816	0	856	2	0
1	E	816	0	856	5	0
1	K	816	0	856	5	0
1	O	825	0	869	6	0
2	B	638	0	676	6	0
2	F	638	0	676	4	0
2	L	638	0	676	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	P	638	0	676	2	0
3	C	810	0	866	3	0
3	G	796	0	848	8	0
3	M	810	0	866	5	0
3	Q	805	0	861	5	0
4	D	788	0	826	5	0
4	H	788	0	826	7	0
4	N	788	0	826	5	0
4	R	788	0	826	6	0
5	I	3462	0	1901	21	0
5	S	3462	0	1901	37	0
6	J	3461	0	1902	12	0
6	T	3461	0	1902	21	0
7	U	560	0	619	36	0
8	G	1	0	0	0	0
8	I	3	0	0	0	0
8	J	1	0	0	0	0
8	Q	3	0	0	0	0
8	S	2	0	0	0	0
8	T	3	0	0	0	0
All	All	26617	0	21111	169	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 4.

All (169) 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)
5:S:-77:DT:H72	7:U:41:VAL:HG23	1.55	0.89
7:U:45:ILE:HD12	7:U:67:LEU:HD21	1.58	0.86
4:R:49:HIS:HB3	4:R:52:THR:HG23	1.59	0.84
1:K:76:GLN:OE1	2:L:23:ARG:NH1	2.14	0.79
5:S:-75:DT:H71	7:U:44:LEU:HD22	1.65	0.76
5:S:-77:DT:C7	7:U:41:VAL:HG23	2.15	0.76
6:T:80:DA:H2''	6:T:81:DG:O5'	1.90	0.71
5:S:-77:DT:H72	7:U:41:VAL:CG2	2.24	0.66
3:Q:81:ARG:NH2	3:Q:107:VAL:O	2.27	0.64
7:U:60:LEU:HD11	7:U:81:ILE:HD12	1.80	0.63
6:T:-62:DC:H2'	6:T:-61:DG:O4'	1.99	0.62
4:D:95:GLN:HE21	4:D:111:VAL:HG13	1.65	0.62
5:S:83:DT:H2''	5:S:84:DG:C8	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:7:DC:H2''	6:T:8:DG:C8	2.36	0.61
7:U:64:LYS:HG2	7:U:81:ILE:HD13	1.83	0.60
6:T:72:DA:OP1	7:U:51:ALA:HB1	2.02	0.59
7:U:64:LYS:CG	7:U:81:ILE:HD13	2.33	0.58
5:S:-47:DT:H2''	5:S:-46:DC:C6	2.39	0.58
5:S:49:DC:H2'	5:S:50:DG:C8	2.39	0.58
6:T:54:DT:H2''	6:T:55:DC:C6	2.39	0.57
2:L:68:ASP:OD2	2:L:92:ARG:NH1	2.37	0.57
5:I:82:DC:H2'	5:I:83:DT:C6	2.40	0.57
5:I:80:DA:H2''	5:I:81:DG:OP2	2.04	0.57
5:S:-77:DT:C7	7:U:41:VAL:CG2	2.82	0.56
6:J:75:DA:H2''	6:J:76:DA:OP2	2.05	0.56
7:U:94:LEU:HD23	7:U:106:PHE:CB	2.36	0.56
7:U:94:LEU:HD23	7:U:106:PHE:HB2	1.88	0.56
7:U:63:LEU:HD12	7:U:106:PHE:CZ	2.42	0.55
1:A:96:CYS:SG	2:B:62:LEU:HD21	2.47	0.55
7:U:45:ILE:HD11	7:U:49:VAL:HG23	1.87	0.54
7:U:88:LEU:HD12	7:U:89:VAL:HG23	1.89	0.54
5:I:11:DC:H2'	5:I:12:DG:C8	2.41	0.54
5:S:-16:DT:H2''	5:S:-15:DA:C8	2.42	0.54
5:S:75:DA:H2''	5:S:76:DA:OP2	2.06	0.54
7:U:45:ILE:HD12	7:U:67:LEU:CD2	2.35	0.54
5:I:-82:DG:H8	5:S:86:DA:C2	2.25	0.53
2:F:68:ASP:OD2	2:F:92:ARG:NH1	2.40	0.53
5:I:51:DG:H2''	5:I:52:DC:H5''	1.91	0.53
5:I:85:DC:N4	6:T:83:DT:O4	2.41	0.53
5:I:49:DC:H2'	5:I:50:DG:C8	2.44	0.53
2:B:24:ASP:HB3	2:B:26:ILE:HG22	1.90	0.53
4:H:49:HIS:HB3	4:H:52:THR:CG2	2.38	0.52
4:R:49:HIS:HB3	4:R:52:THR:CG2	2.35	0.52
4:H:33:ARG:O	4:H:33:ARG:HG3	2.09	0.52
7:U:97:THR:HG23	7:U:107:LYS:HG2	1.92	0.52
7:U:40:PRO:O	7:U:41:VAL:HG22	2.10	0.52
6:T:49:DC:H2'	6:T:50:DG:C8	2.45	0.51
5:S:35:DT:H1'	5:S:36:DA:C5	2.46	0.51
7:U:51:ALA:HA	7:U:108:LEU:HD21	1.92	0.51
1:A:119:ILE:HD12	2:B:50:ILE:HD13	1.93	0.51
2:B:68:ASP:OD2	2:B:92:ARG:NH1	2.44	0.51
4:H:49:HIS:HB3	4:H:52:THR:HG22	1.93	0.51
4:R:65:PHE:CE1	4:R:69:ILE:CD1	2.95	0.50
6:J:7:DC:H2''	6:J:8:DG:C8	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:82:DC:N4	6:T:86:DA:N6	2.60	0.49
5:S:54:DT:H2''	5:S:55:DC:C6	2.46	0.49
7:U:94:LEU:HD23	7:U:106:PHE:CG	2.47	0.49
5:S:20:DG:H2''	5:S:21:DG:O5'	2.12	0.49
7:U:39:PRO:HG3	7:U:67:LEU:HA	1.95	0.49
6:J:66:DT:H1'	6:J:67:DT:H5'	1.94	0.49
5:S:-73:DT:H1'	5:S:-72:DT:H5'	1.94	0.49
5:I:80:DA:C2'	5:I:81:DG:OP2	2.61	0.49
5:I:-81:DC:H2'	5:I:-80:DT:C6	2.48	0.49
6:T:-7:DG:OP2	7:U:107:LYS:NZ	2.45	0.49
4:D:30:LYS:NZ	5:I:-46:DC:OP1	2.45	0.48
1:K:119:ILE:HD12	2:L:50:ILE:HD13	1.95	0.48
4:N:65:PHE:CE1	4:N:69:ILE:CD1	2.95	0.48
3:G:17:ARG:NH2	3:G:28:GLY:HA2	2.27	0.48
6:T:11:DC:H2'	6:T:12:DG:C8	2.49	0.48
5:S:-78:DT:H2''	5:S:-77:DT:C5	2.47	0.48
3:G:44:GLY:HA2	5:I:38:DG:H5''	1.95	0.48
5:S:48:DG:H1'	5:S:49:DC:C6	2.49	0.47
5:S:-76:DT:H4'	5:S:-75:DT:OP1	2.14	0.47
3:C:100:VAL:HG11	2:F:98:TYR:CE2	2.50	0.47
1:E:56:LYS:NZ	5:I:-64:DC:OP1	2.48	0.47
7:U:51:ALA:HB2	7:U:108:LEU:CD2	2.44	0.46
5:S:-2:DG:H1'	5:S:-1:DA:C8	2.51	0.46
3:G:16:THR:HG23	3:G:19:SER:CB	2.45	0.46
2:B:24:ASP:CB	2:B:26:ILE:HG22	2.45	0.46
3:M:115:LEU:HD11	1:O:108:ASN:HD21	1.81	0.46
5:S:-82:DG:H2''	5:S:-81:DC:O4'	2.14	0.46
1:E:131:ARG:HD3	1:E:133:GLU:OE2	2.16	0.46
7:U:95:VAL:HB	7:U:107:LYS:HE2	1.98	0.46
3:G:24:GLN:HG3	4:H:43:LYS:HB3	1.98	0.46
2:L:98:TYR:CE2	3:Q:100:VAL:HG11	2.51	0.45
4:R:124:ALA:O	4:R:125:LYS:HB2	2.17	0.45
5:S:82:DC:H2''	5:S:83:DT:H5'	1.98	0.45
3:C:115:LEU:HD11	1:E:108:ASN:HD21	1.81	0.45
7:U:48:ALA:O	7:U:52:SER:N	2.49	0.45
5:I:81:DG:C5	5:I:82:DC:C4	3.04	0.45
5:I:-16:DT:H2''	5:I:-15:DA:C8	2.52	0.45
2:L:66:ILE:O	2:L:67:ARG:C	2.60	0.45
6:T:73:DA:OP1	7:U:52:SER:OG	2.20	0.45
4:D:118:VAL:O	4:D:122:THR:OG1	2.35	0.45
5:I:55:DC:H2''	5:I:56:DG:N7	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:-62:DC:H2'	5:S:-61:DG:C8	2.52	0.45
7:U:63:LEU:HD12	7:U:106:PHE:CE2	2.52	0.45
6:J:-6:DT:H2''	6:J:-5:DA:C8	2.52	0.44
4:D:76:GLU:OE2	4:D:79:ARG:NH1	2.50	0.44
1:O:42:ARG:NH2	6:T:71:DA:OP2	2.50	0.44
5:S:-75:DT:C7	7:U:44:LEU:HD22	2.43	0.44
6:T:76:DA:C2	6:T:77:DA:C4	3.05	0.44
2:P:60:VAL:O	2:P:64:ASN:ND2	2.44	0.44
4:R:95:GLN:NE2	4:R:111:VAL:HG13	2.32	0.44
6:J:74:DA:H2''	6:J:75:DA:C8	2.52	0.44
6:T:-47:DT:H2''	6:T:-46:DC:C6	2.53	0.44
5:I:-65:DT:H1'	5:I:-64:DC:H5''	2.00	0.44
6:J:83:DT:H2''	6:J:84:DG:O4'	2.18	0.44
3:G:102:ILE:HG23	4:H:61:ILE:HD13	1.99	0.44
1:O:36:LYS:N	5:S:-68:DC:OP1	2.51	0.44
6:J:-7:DG:H2''	6:J:-6:DT:H71	2.00	0.43
3:M:90:ASP:O	3:M:91:GLU:C	2.60	0.43
5:S:83:DT:H4'	5:S:84:DG:OP1	2.18	0.43
2:L:88:TYR:CE1	4:N:83:TYR:CE1	3.06	0.43
3:M:79:ILE:HG12	3:M:82:HIS:CE1	2.53	0.43
1:E:69:ARG:NH2	5:I:17:DA:OP2	2.52	0.43
5:S:80:DA:C6	5:S:81:DG:C6	3.07	0.43
4:D:95:GLN:NE2	4:D:111:VAL:HG13	2.33	0.43
3:G:50:TYR:O	3:G:54:VAL:HG23	2.19	0.43
2:L:84:MET:HE3	2:L:88:TYR:CZ	2.54	0.43
7:U:82:LYS:HE2	7:U:102:ALA:HB1	2.00	0.43
1:E:91:ALA:HB2	2:F:100:PHE:CD2	2.54	0.43
1:K:119:ILE:HG13	2:L:50:ILE:CD1	2.49	0.43
5:S:-56:DC:O5'	5:S:-56:DC:H2'	2.19	0.43
6:T:-46:DC:H2''	6:T:-45:DA:C8	2.54	0.42
7:U:45:ILE:HA	7:U:48:ALA:HB3	2.01	0.42
2:L:30:THR:HB	2:L:32:PRO:HD2	2.02	0.42
6:T:-74:DT:C6	6:T:-73:DT:H72	2.54	0.42
6:T:-46:DC:H2''	6:T:-45:DA:N7	2.33	0.42
3:M:31:HIS:CD2	3:M:48:PRO:HG3	2.54	0.42
1:O:62:ILE:HG21	2:P:29:ILE:HD12	2.02	0.42
5:S:-78:DT:H2''	5:S:-77:DT:C7	2.49	0.42
1:K:42:ARG:HD3	5:S:71:DA:OP2	2.19	0.42
3:Q:31:HIS:CD2	3:Q:48:PRO:HG3	2.54	0.42
7:U:40:PRO:CG	7:U:44:LEU:HD11	2.49	0.42
4:N:86:ARG:NH2	5:S:-33:DA:OP2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:63:LEU:HD22	7:U:67:LEU:HD11	2.01	0.42
6:J:-80:DT:H2''	6:J:-79:DT:OP2	2.19	0.42
3:Q:80:PRO:HG3	4:R:61:ILE:HD12	2.01	0.42
7:U:40:PRO:HB2	7:U:44:LEU:HD11	2.01	0.42
5:S:-3:DG:C6	5:S:-2:DG:C6	3.07	0.42
6:T:15:DT:H2''	6:T:16:DA:C8	2.55	0.42
5:S:82:DC:C2'	5:S:83:DT:O5'	2.68	0.42
2:F:88:TYR:CE1	4:H:83:TYR:CE1	3.08	0.42
3:Q:18:SER:OG	3:Q:25:PHE:O	2.38	0.42
5:S:22:DT:H2'	5:S:23:DG:C8	2.54	0.42
1:O:92:LEU:HD23	1:O:92:LEU:HA	1.92	0.42
6:T:-54:DA:C6	6:T:-53:DG:C6	3.07	0.42
5:I:28:DA:H2''	5:I:29:DG:H5'	2.02	0.41
6:J:7:DC:C2'	6:J:8:DG:C8	3.02	0.41
6:J:-1:DA:H1'	6:J:0:DT:H5'	2.02	0.41
3:C:80:PRO:HB2	3:C:104:GLN:O	2.20	0.41
5:S:-26:DT:C4	5:S:-25:DA:C6	3.08	0.41
6:J:-79:DT:H2''	6:J:-78:DT:H5''	2.02	0.41
3:M:64:GLU:OE1	4:N:48:VAL:HG13	2.19	0.41
4:N:42:TYR:OH	5:S:-53:DG:OP2	2.30	0.41
1:O:83:ARG:HG2	6:T:-24:DG:O3'	2.21	0.41
2:B:84:MET:SD	2:B:101:GLY:O	2.79	0.41
5:I:86:DA:C2'	5:S:-82:DG:H5'	2.51	0.41
6:J:-77:DT:H2''	6:J:-76:DT:H6	1.86	0.41
6:T:-34:DG:H1'	6:T:-33:DA:C8	2.56	0.41
7:U:93:THR:HG22	7:U:94:LEU:HD13	2.03	0.41
3:G:16:THR:HG23	3:G:19:SER:HB2	2.03	0.41
7:U:82:LYS:CE	7:U:102:ALA:HB1	2.51	0.41
3:G:29:ARG:NH2	4:H:36:SER:O	2.55	0.40
2:L:31:LYS:HG3	2:L:51:TYR:CE1	2.56	0.40
7:U:81:ILE:HG13	7:U:82:LYS:N	2.36	0.40
5:I:28:DA:H1'	5:I:29:DG:H5'	2.04	0.40
1:K:83:ARG:HA	5:S:-23:DC:OP1	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	97/138 (70%)	94 (97%)	3 (3%)	0	100	100
1	E	97/138 (70%)	95 (98%)	2 (2%)	0	100	100
1	K	97/138 (70%)	96 (99%)	1 (1%)	0	100	100
1	O	98/138 (71%)	96 (98%)	2 (2%)	0	100	100
2	B	78/105 (74%)	76 (97%)	2 (3%)	0	100	100
2	F	78/105 (74%)	75 (96%)	3 (4%)	0	100	100
2	L	78/105 (74%)	73 (94%)	5 (6%)	0	100	100
2	P	78/105 (74%)	76 (97%)	2 (3%)	0	100	100
3	C	103/132 (78%)	97 (94%)	6 (6%)	0	100	100
3	G	101/132 (76%)	98 (97%)	3 (3%)	0	100	100
3	M	103/132 (78%)	97 (94%)	6 (6%)	0	100	100
3	Q	102/132 (77%)	97 (95%)	5 (5%)	0	100	100
4	D	98/128 (77%)	93 (95%)	3 (3%)	2 (2%)	6	31
4	H	98/128 (77%)	88 (90%)	8 (8%)	2 (2%)	6	31
4	N	98/128 (77%)	92 (94%)	5 (5%)	1 (1%)	12	45
4	R	98/128 (77%)	92 (94%)	3 (3%)	3 (3%)	3	22
7	U	76/222 (34%)	62 (82%)	12 (16%)	2 (3%)	4	26
All	All	1578/2234 (71%)	1497 (95%)	71 (4%)	10 (1%)	21	56

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	H	124	ALA
7	U	41	VAL
7	U	77	ASN
4	R	104	GLY
4	D	29	ARG

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Mol	Chain	Res	Type
4	N	104	GLY
4	R	33	ARG
4	D	104	GLY
4	R	54	ILE
4	H	54	ILE

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	86/112 (77%)	84 (98%)	2 (2%)	44	70
1	E	86/112 (77%)	82 (95%)	4 (5%)	23	57
1	K	86/112 (77%)	82 (95%)	4 (5%)	23	57
1	O	87/112 (78%)	81 (93%)	6 (7%)	14	45
2	B	65/80 (81%)	65 (100%)	0	100	100
2	F	65/80 (81%)	64 (98%)	1 (2%)	57	76
2	L	65/80 (81%)	63 (97%)	2 (3%)	35	66
2	P	65/80 (81%)	64 (98%)	1 (2%)	57	76
3	C	83/101 (82%)	79 (95%)	4 (5%)	23	56
3	G	82/101 (81%)	76 (93%)	6 (7%)	13	43
3	M	83/101 (82%)	77 (93%)	6 (7%)	13	44
3	Q	83/101 (82%)	76 (92%)	7 (8%)	10	38
4	D	85/106 (80%)	77 (91%)	8 (9%)	8	33
4	H	85/106 (80%)	77 (91%)	8 (9%)	8	33
4	N	85/106 (80%)	78 (92%)	7 (8%)	10	39
4	R	85/106 (80%)	77 (91%)	8 (9%)	8	33
7	U	59/159 (37%)	42 (71%)	17 (29%)	0	1
All	All	1335/1755 (76%)	1244 (93%)	91 (7%)	14	46

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LYS
1	A	129	ARG
3	C	15	LYS
3	C	24	GLN
3	C	74	LYS
3	C	81	ARG
4	D	32	SER
4	D	34	LYS
4	D	36	SER
4	D	63	ASN
4	D	85	LYS
4	D	101	LEU
4	D	106	LEU
4	D	122	THR
1	E	37	LYS
1	E	48	LEU
1	E	129	ARG
1	E	134	ARG
2	F	27	GLN
3	G	20	ARG
3	G	24	GLN
3	G	29	ARG
3	G	71	ARG
3	G	74	LYS
3	G	81	ARG
4	H	27	LYS
4	H	46	LYS
4	H	52	THR
4	H	80	LEU
4	H	85	LYS
4	H	105	GLU
4	H	106	LEU
4	H	125	LYS
1	K	37	LYS
1	K	63	ARG
1	K	86	SER
1	K	129	ARG
2	L	24	ASP
2	L	50	ILE
3	M	15	LYS
3	M	24	GLN
3	M	36	LYS
3	M	73	ASN

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Mol	Chain	Res	Type
3	M	74	LYS
3	M	81	ARG
4	N	29	ARG
4	N	32	SER
4	N	36	SER
4	N	63	ASN
4	N	85	LYS
4	N	101	LEU
4	N	125	LYS
1	O	37	LYS
1	O	48	LEU
1	O	76	GLN
1	O	83	ARG
1	O	129	ARG
1	O	134	ARG
2	P	27	GLN
3	Q	24	GLN
3	Q	29	ARG
3	Q	42	ARG
3	Q	64	GLU
3	Q	74	LYS
3	Q	81	ARG
3	Q	118	LYS
4	R	28	LYS
4	R	32	SER
4	R	36	SER
4	R	46	LYS
4	R	52	THR
4	R	56	SER
4	R	80	LEU
4	R	85	LYS
7	U	34	ARG
7	U	44	LEU
7	U	45	ILE
7	U	49	VAL
7	U	58	VAL
7	U	63	LEU
7	U	64	LYS
7	U	75	GLU
7	U	77	ASN
7	U	78	ASN
7	U	85	LEU

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Mol	Chain	Res	Type
7	U	87	SER
7	U	90	SER
7	U	96	GLN
7	U	100	THR
7	U	108	LEU
7	U	109	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	76	GLN
3	C	24	GLN
3	C	31	HIS
4	D	47	GLN
4	D	84	ASN
4	D	95	GLN
1	E	39	HIS
1	E	76	GLN
1	E	85	GLN
1	E	108	ASN
3	G	94	ASN
4	H	67	ASN
3	M	24	GLN
3	M	31	HIS
3	M	104	GLN
4	N	47	GLN
4	N	95	GLN
1	O	39	HIS
1	O	108	ASN
3	Q	24	GLN
3	Q	31	HIS
3	Q	68	ASN
4	R	95	GLN
7	U	109	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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(Å²)	Q<0.9	
1	A	99/138 (71%)	-0.65	0	100	100	99, 119, 148, 165	0
1	E	99/138 (71%)	-0.70	0	100	100	95, 112, 148, 161	0
1	K	99/138 (71%)	-0.74	0	100	100	57, 80, 117, 136	0
1	O	100/138 (72%)	-0.75	0	100	100	57, 73, 109, 141	0
2	B	80/105 (76%)	-0.63	0	100	100	101, 118, 139, 168	0
2	F	80/105 (76%)	-0.62	0	100	100	96, 110, 132, 154	0
2	L	80/105 (76%)	-0.81	0	100	100	61, 79, 106, 163	0
2	P	80/105 (76%)	-0.86	0	100	100	57, 73, 95, 133	0
3	C	105/132 (79%)	-0.54	0	100	100	98, 117, 148, 156	0
3	G	103/132 (78%)	-0.56	0	100	100	96, 120, 147, 162	0
3	M	105/132 (79%)	-0.66	0	100	100	62, 78, 107, 133	0
3	Q	104/132 (78%)	-0.65	0	100	100	64, 80, 114, 140	0
4	D	100/128 (78%)	-0.50	0	100	100	99, 120, 165, 179	0
4	H	100/128 (78%)	-0.51	0	100	100	101, 121, 171, 192	0
4	N	100/128 (78%)	-0.67	0	100	100	67, 83, 143, 171	0
4	R	100/128 (78%)	-0.62	0	100	100	67, 82, 157, 179	0
5	I	169/169 (100%)	-0.71	0	100	100	122, 163, 219, 253	0
5	S	169/169 (100%)	-0.70	0	100	100	83, 125, 211, 260	0
6	J	169/169 (100%)	-0.71	0	100	100	122, 161, 209, 273	0
6	T	169/169 (100%)	-0.75	0	100	100	84, 124, 200, 250	0
7	U	78/222 (35%)	0.50	2 (2%)	57	37	151, 170, 223, 263	0
All	All	2288/2910 (78%)	-0.63	2 (0%)	92	89	57, 114, 187, 273	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
7	U	47	LYS	3.5
7	U	49	VAL	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	CA	I	103	1/1	0.88	0.07	120,120,120,120	0
8	CA	G	201	1/1	0.89	0.07	129,129,129,129	0
8	CA	T	103	1/1	0.89	0.07	124,124,124,124	0
8	CA	S	102	1/1	0.90	0.05	112,112,112,112	0
8	CA	Q	203	1/1	0.91	0.06	115,115,115,115	0
8	CA	T	102	1/1	0.92	0.12	126,126,126,126	0
8	CA	T	101	1/1	0.95	0.11	92,92,92,92	0
8	CA	S	101	1/1	0.95	0.05	107,107,107,107	0
8	CA	I	102	1/1	0.95	0.04	113,113,113,113	0
8	CA	I	101	1/1	0.96	0.10	112,112,112,112	0
8	CA	J	101	1/1	0.96	0.08	119,119,119,119	0
8	CA	Q	202	1/1	0.96	0.04	116,116,116,116	0
8	CA	Q	201	1/1	0.97	0.03	96,96,96,96	0

6.5 Other polymers [i](#)

There are no such residues in this entry.