Full wwPDB X-ray Structure Validation Report (i)

Oct 30, 2023 – 08:42 PM JST

PDB ID	:	4Z5T
Title	:	The nucleosome containing human H3.5
Authors	:	Urahama, T.; Harada, A.; Maehara, K.; Horikoshi, N.; Sato, K.; Sato, Y.; Shiraishi, K.; Sugino, N.; Osakabe, A.; Tachiwana, H.; Kagawa, W.; Kimura, H.; Ohkawa, Y.; Kurumizaka, H.
Deposited on	:	2015-04-03
Resolution	:	2.80 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 (i) 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 (1)) were used in the production of this report:

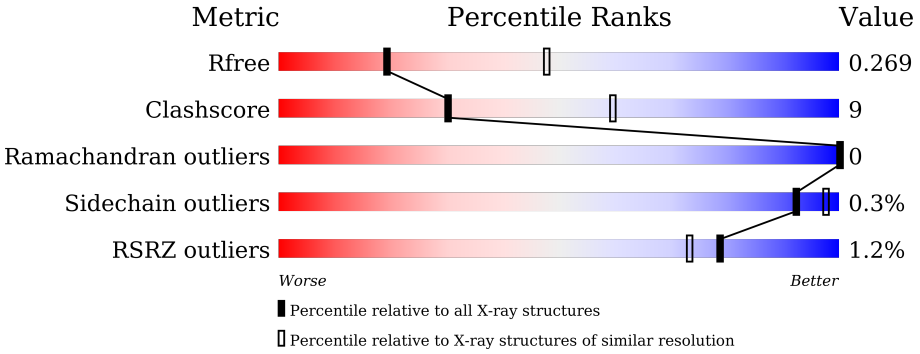
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	.% 56% 14% 30%
1	E	138	.% 56% 14% 30%
2	B	106	68% 10% 22%
2	F	106	2% 56% 19% 25%
3	C	133	.% 67% 12% 21%
3	G	133	.% 64% 14% • 22%

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Mol	Chain	Length	Quality of chain
4	D	129	% 60% 11% 29%
4	H	129	64% 8% 29%
5	I	146	% 40% 60%
5	J	146	% 49% 51%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 11899 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 H3.3C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	97	792	497	155	138	2	0	0	0
1	E	96	785	492	154	137	2	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q6NXT2
A	-2	SER	-	expression tag	UNP Q6NXT2
A	-1	HIS	-	expression tag	UNP Q6NXT2
E	-3	GLY	-	expression tag	UNP Q6NXT2
E	-2	SER	-	expression tag	UNP Q6NXT2
E	-1	HIS	-	expression tag	UNP Q6NXT2

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	83	662	418	129	114	1	0	0	0
2	F	79	627	395	121	110	1	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P62805
B	-2	SER	-	expression tag	UNP P62805
B	-1	HIS	-	expression tag	UNP P62805
F	-3	GLY	-	expression tag	UNP P62805
F	-2	SER	-	expression tag	UNP P62805
F	-1	HIS	-	expression tag	UNP P62805

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	105	810	511	158	141	0	0	0
3	G	104	805	508	157	140	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLY	-	expression tag	UNP P04908
C	-2	SER	-	expression tag	UNP P04908
C	-1	HIS	-	expression tag	UNP P04908
G	-3	GLY	-	expression tag	UNP P04908
G	-2	SER	-	expression tag	UNP P04908
G	-1	HIS	-	expression tag	UNP P04908

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	92	719	453	129	135	2	0	0	0
4	H	92	719	453	129	135	2	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-3	GLY	-	expression tag	UNP P06899
D	-2	SER	-	expression tag	UNP P06899
D	-1	HIS	-	expression tag	UNP P06899
H	-3	GLY	-	expression tag	UNP P06899
H	-2	SER	-	expression tag	UNP P06899
H	-1	HIS	-	expression tag	UNP P06899

- Molecule 5 is a DNA chain called DNA (146-MER).

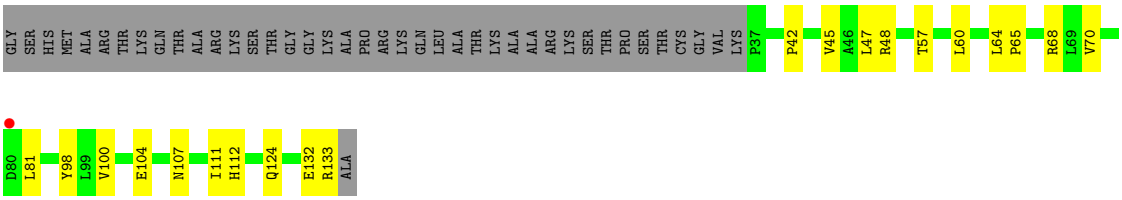
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
5	I	146	2990	1431	540	874	145	0	0	0
5	J	146	2990	1431	540	874	145	0	0	0

3 Residue-property plots (i)

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.

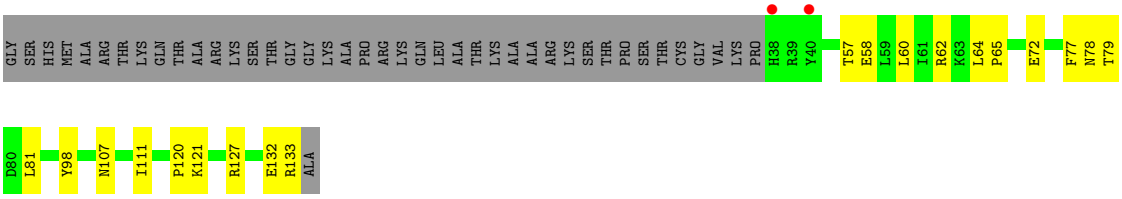
- Molecule 1: Histone H3.3C

Chain A:



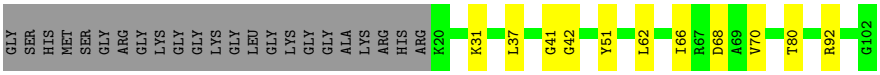
- Molecule 1: Histone H3.3C

Chain E:



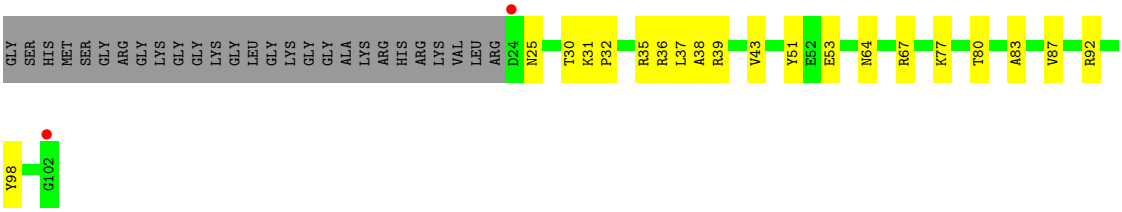
- Molecule 2: Histone H4

Chain B:



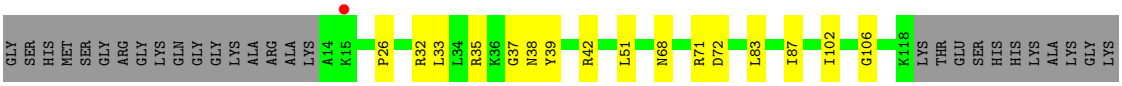
- Molecule 2: Histone H4

Chain F:



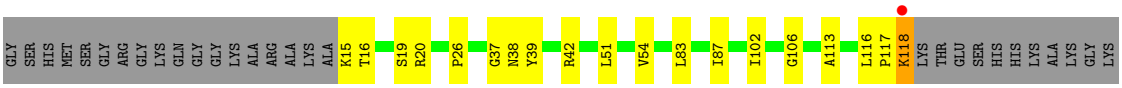
• Molecule 3: Histone H2A type 1-B/E

Chain C:



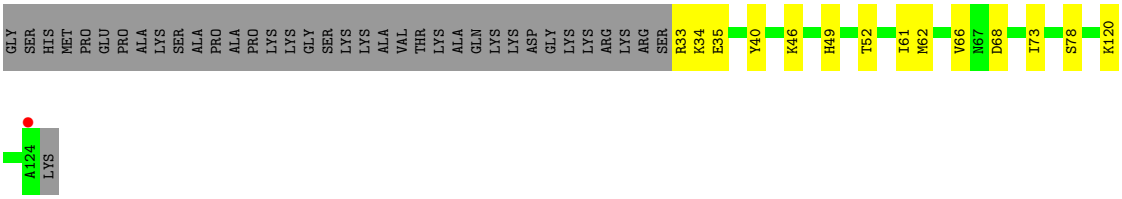
• Molecule 3: Histone H2A type 1-B/E

Chain G:



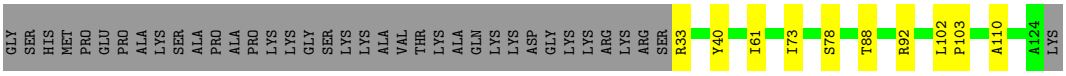
• Molecule 4: Histone H2B type 1-J

Chain D:



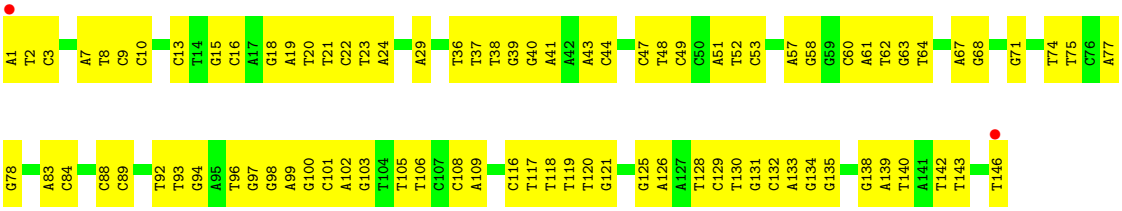
• Molecule 4: Histone H2B type 1-J

Chain H:



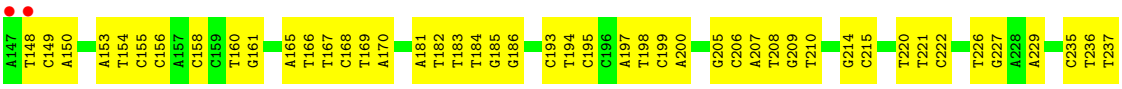
• Molecule 5: DNA (146-MER)

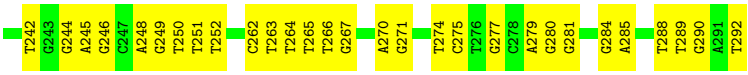
Chain I:



• Molecule 5: DNA (146-MER)

Chain J:




T242
G243
G244
A245
G246
C247
A248
G249
T250
T251
T252
C262
T263
T264
T265
T266
G267
A270
G271
T274
C275
T276
G277
C278
A279
G280
G281
G284
A285
T288
T289
G290
A291
T292

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	104.89Å 109.13Å 174.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.81 – 2.80 47.27 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.8 (37.81-2.80) 99.2 (47.27-2.78)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.18 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.228 , 0.269 0.228 , 0.269	Depositor DCC
R_{free} test set	2570 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	60.4	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 61.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11899	wwPDB-VP
Average B, all atoms (Å^2)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.32	0/803	0.52	0/1079
1	E	0.31	0/795	0.50	0/1068
2	B	0.28	0/669	0.48	0/894
2	F	0.26	0/634	0.44	0/848
3	C	0.26	0/820	0.47	0/1107
3	G	0.29	0/815	0.52	0/1100
4	D	0.29	0/730	0.49	0/982
4	H	0.30	0/730	0.47	0/982
5	I	0.52	0/3354	0.94	0/5175
5	J	0.55	0/3354	0.96	0/5175
All	All	0.44	0/12704	0.78	0/18410

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	792	0	828	17	0
1	E	785	0	820	17	1
2	B	662	0	709	11	0
2	F	627	0	663	15	0
3	C	810	0	866	15	0
3	G	805	0	861	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	719	0	740	12	0
4	H	719	0	740	11	0
5	I	2990	0	1652	69	0
5	J	2990	0	1652	69	1
All	All	11899	0	9531	199	1

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 9.

All (199) 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:E:62:ARG:HH21	5:J:207:DA:H5'	1.51	0.73
5:I:23:DT:H2"	5:I:24:DA:H5"	1.70	0.71
5:I:8:DT:H2"	5:I:9:DC:H5"	1.72	0.70
5:J:182:DT:H2"	5:J:183:DT:H5"	1.74	0.70
5:J:288:DT:H1'	5:J:289:DT:H5'	1.74	0.69
5:J:155:DC:H2"	5:J:156:DC:H5'	1.75	0.69
5:I:40:DG:H2"	5:I:41:DA:C8	2.29	0.68
5:J:279:DA:H2"	5:J:280:DG:H5"	1.74	0.68
5:I:9:DC:H42	5:J:284:DG:H1	1.42	0.67
5:J:149:DC:H1'	5:J:150:DA:H5'	1.77	0.67
5:I:9:DC:H2"	5:I:10:DC:H5'	1.77	0.66
5:J:220:DT:H1'	5:J:221:DT:H5'	1.77	0.66
3:G:37:GLY:HA3	3:G:39:TYR:CE2	2.32	0.64
5:J:205:DG:H2"	5:J:206:DC:H5"	1.79	0.63
5:I:36:DT:H2"	5:I:37:DT:H5"	1.79	0.63
5:J:274:DT:H2"	5:J:275:DC:H5"	1.81	0.63
5:I:98:DG:H1	5:J:195:DC:H42	1.46	0.63
5:I:74:DT:H1'	5:I:75:DT:H5'	1.82	0.62
5:I:142:DT:H1'	5:I:143:DT:H5'	1.82	0.61
3:G:15:LYS:O	3:G:20:ARG:NH1	2.34	0.61
5:I:67:DA:H2"	5:I:68:DG:C8	2.36	0.61
1:E:60:LEU:HD12	2:F:37:LEU:HD23	1.82	0.60
1:E:120:PRO:HB3	2:F:53:GLU:HG3	1.82	0.60
5:I:105:DT:H1'	5:I:106:DT:H5'	1.84	0.60
5:I:125:DG:H2"	5:I:126:DA:H8	1.66	0.60
5:J:194:DT:H2"	5:J:195:DC:H5'	1.84	0.60
1:E:62:ARG:NH2	5:J:206:DC:O3'	2.35	0.60
3:C:42:ARG:CZ	5:I:38:DT:H5"	2.32	0.59
5:J:153:DA:H1'	5:J:154:DT:H5'	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:33:ARG:HG2	4:D:34:LYS:HG3	1.83	0.58
5:J:245:DA:H2"	5:J:246:DG:C8	2.38	0.58
5:J:198:DT:H2"	5:J:199:DC:H5'	1.86	0.58
5:J:270:DA:H2"	5:J:271:DG:O5'	2.04	0.58
5:I:133:DA:H2"	5:I:134:DG:H5"	1.85	0.57
1:A:107:ASN:O	1:A:111:ILE:HG12	2.04	0.57
3:G:113:ALA:O	3:G:118:LYS:NZ	2.38	0.57
5:J:197:DA:H2"	5:J:198:DT:H5"	1.87	0.57
5:I:132:DC:H2"	5:I:133:DA:C8	2.39	0.57
5:J:266:DT:H2"	5:J:267:DG:C8	2.40	0.57
5:J:148:DT:H2"	5:J:149:DC:H5'	1.87	0.56
5:J:209:DG:H1'	5:J:210:DT:H5'	1.87	0.56
4:H:33:ARG:HG2	4:H:33:ARG:NH1	2.19	0.56
1:A:60:LEU:HD12	2:B:37:LEU:HD23	1.88	0.56
1:E:127:ARG:HH21	1:E:133:ARG:HD2	1.71	0.55
5:J:165:DA:H2"	5:J:166:DT:H5'	1.89	0.55
5:J:207:DA:H2"	5:J:208:DT:H5'	1.89	0.55
3:C:32:ARG:HH21	3:C:33:LEU:HD21	1.72	0.55
5:I:52:DT:H2"	5:I:53:DC:H5'	1.88	0.55
3:G:102:ILE:HG23	4:H:61:ILE:HD13	1.90	0.54
5:J:226:DT:H2"	5:J:227:DG:C8	2.43	0.54
5:I:47:DC:H2"	5:I:48:DT:C6	2.42	0.54
5:I:13:DC:H42	5:J:280:DG:H1	1.55	0.53
1:A:100:VAL:O	1:A:104:GLU:HG3	2.08	0.53
1:E:58:GLU:OE2	1:E:58:GLU:N	2.35	0.53
2:B:31:LYS:HG3	2:B:51:TYR:CE1	2.44	0.53
1:E:62:ARG:HB2	1:E:65:PRO:HD2	1.91	0.52
5:I:102:DA:H2"	5:I:103:DG:H8	1.74	0.52
5:I:7:DA:H1'	5:I:8:DT:H5'	1.92	0.52
5:I:128:DT:H2"	5:I:129:DC:H5"	1.92	0.52
5:I:120:DT:H2"	5:I:121:DG:C8	2.45	0.52
5:J:248:DA:H2"	5:J:249:DG:H8	1.75	0.52
3:G:116:LEU:HB2	3:G:118:LYS:NZ	2.25	0.51
4:H:33:ARG:HG2	4:H:33:ARG:HH11	1.74	0.51
5:I:38:DT:H2"	5:I:39:DG:N7	2.25	0.51
5:I:108:DC:H2"	5:I:109:DA:N7	2.25	0.51
5:I:116:DC:H1'	5:I:117:DT:H5'	1.92	0.51
5:J:183:DT:H2"	5:J:184:DT:H5'	1.93	0.51
1:A:47:LEU:HD13	3:G:117:PRO:HD3	1.92	0.51
5:I:130:DT:H2"	5:I:131:DG:C8	2.46	0.51
5:J:251:DT:H1'	5:J:252:DT:H5'	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:125:DG:H2"	5:I:126:DA:C8	2.45	0.50
2:B:31:LYS:HG3	2:B:51:TYR:CZ	2.46	0.50
5:J:277:DG:H5"	5:J:277:DG:H8	1.75	0.50
5:I:9:DC:N4	5:J:284:DG:H1	2.08	0.50
1:A:42:PRO:HA	5:J:229:DA:H5"	1.94	0.50
2:F:77:LYS:HE3	4:H:92:ARG:CZ	2.42	0.50
5:I:1:DA:H61	5:J:292:DT:H3	1.58	0.50
5:I:138:DG:N2	5:J:156:DC:O2	2.45	0.50
3:G:37:GLY:O	3:G:38:ASN:HB2	2.12	0.49
5:I:99:DA:H2"	5:I:100:DG:C8	2.47	0.49
1:E:127:ARG:NH2	1:E:133:ARG:HD2	2.27	0.49
3:C:71:ARG:HG3	3:C:72:ASP:N	2.26	0.49
5:I:51:DA:H2"	5:I:52:DT:H5"	1.94	0.49
5:I:135:DG:H1	5:J:158:DC:H42	1.60	0.49
5:J:155:DC:C2'	5:J:156:DC:H5'	2.43	0.49
5:I:93:DT:H1'	5:I:94:DG:H5'	1.94	0.49
3:C:83:LEU:O	3:C:87:ILE:HG12	2.13	0.48
1:E:98:TYR:OH	1:E:132:GLU:OE1	2.29	0.48
5:J:167:DT:H1'	5:J:168:DC:H5'	1.95	0.48
2:F:35:ARG:O	2:F:39:ARG:HG2	2.14	0.48
5:J:264:DT:H1'	5:J:265:DT:H5'	1.96	0.48
3:C:102:ILE:HG23	4:D:61:ILE:HD13	1.96	0.48
3:G:39:TYR:HB3	4:H:78:SER:HB2	1.96	0.48
5:J:244:DG:H1'	5:J:245:DA:H5'	1.96	0.48
2:B:92:ARG:NH1	2:B:92:ARG:HB3	2.29	0.48
5:I:100:DG:H1	5:J:193:DC:H42	1.62	0.48
3:G:42:ARG:HD2	5:J:185:DG:H5"	1.95	0.48
5:J:235:DC:H2"	5:J:236:DT:C5	2.49	0.48
1:A:45:VAL:HG22	1:A:48:ARG:HH21	1.79	0.47
5:I:108:DC:H2"	5:I:109:DA:C8	2.48	0.47
5:J:160:DT:H2"	5:J:161:DG:C8	2.49	0.47
5:I:88:DC:H2"	5:I:89:DC:C6	2.49	0.47
2:F:38:ALA:HB1	2:F:43:VAL:HB	1.95	0.47
2:F:64:ASN:OD1	2:F:67:ARG:NH2	2.47	0.47
1:A:68:ARG:HH22	5:J:237:DT:P	2.38	0.47
5:I:60:DC:H2"	5:I:61:DA:C8	2.50	0.47
5:I:47:DC:H42	5:J:246:DG:H1	1.62	0.47
2:B:68:ASP:OD2	2:B:92:ARG:NH1	2.47	0.47
1:E:107:ASN:O	1:E:111:ILE:HG12	2.14	0.47
2:F:31:LYS:HG3	2:F:51:TYR:CE1	2.50	0.47
5:I:92:DT:H1'	5:I:93:DT:H5'	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:J:262:DC:H1'	5:J:263:DT:H5'	1.97	0.47
3:C:68:ASN:HA	3:C:71:ARG:HG2	1.97	0.47
1:E:72:GLU:HB2	2:F:25:ASN:HD21	1.80	0.46
5:I:43:DA:H1'	5:I:44:DC:H5"	1.97	0.46
1:E:60:LEU:O	2:F:36:ARG:NH2	2.33	0.46
5:J:199:DC:H1'	5:J:200:DA:H5'	1.97	0.46
1:A:57:THR:HG22	3:G:106:GLY:HA3	1.98	0.46
2:B:80:THR:N	5:J:248:DA:OP1	2.49	0.45
2:F:80:THR:N	5:I:101:DC:OP1	2.50	0.45
5:J:185:DG:H2"	5:J:186:DG:C8	2.51	0.45
5:I:18:DG:H1	5:J:275:DC:H42	1.64	0.45
1:A:64:LEU:HB3	1:A:65:PRO:HD3	1.98	0.45
3:C:106:GLY:HA3	1:E:57:THR:HG22	1.99	0.45
3:C:37:GLY:O	3:C:38:ASN:HB2	2.16	0.45
1:E:64:LEU:HB3	1:E:65:PRO:HD3	1.99	0.45
3:C:32:ARG:HD3	5:I:29:DA:OP2	2.17	0.45
5:I:3:DC:H42	5:J:290:DG:H1	1.65	0.45
5:I:61:DA:H2"	5:I:62:DT:H5'	1.99	0.44
5:J:169:DT:H1'	5:J:170:DA:H5'	1.98	0.44
5:I:71:DG:H1	5:J:222:DC:H42	1.66	0.44
1:A:104:GLU:HG2	2:B:41:GLY:HA2	1.98	0.44
5:I:77:DA:H2"	5:I:78:DG:C8	2.52	0.44
5:J:280:DG:H2"	5:J:281:DG:H5'	1.98	0.44
5:I:2:DT:H2"	5:I:3:DC:H5'	1.98	0.44
5:J:166:DT:H1'	5:J:167:DT:H5'	2.00	0.44
3:G:51:LEU:HD13	4:H:73:ILE:HG21	1.98	0.44
3:G:83:LEU:O	3:G:87:ILE:HG12	2.17	0.44
5:I:20:DT:H1'	5:I:21:DT:H5'	1.98	0.44
5:I:21:DT:H1'	5:I:22:DC:H5'	1.99	0.44
5:I:134:DG:H2"	5:I:135:DG:H5'	2.00	0.44
3:C:26:PRO:HD3	4:D:40:TYR:CD1	2.53	0.43
4:D:49:HIS:HB3	4:D:52:THR:OG1	2.19	0.43
5:J:195:DC:H6	5:J:195:DC:H2'	1.71	0.43
5:I:96:DT:H1'	5:I:97:DG:H5'	1.99	0.43
1:A:98:TYR:OH	1:A:132:GLU:OE1	2.35	0.43
3:C:51:LEU:HD13	4:D:73:ILE:HG21	2.01	0.43
1:A:124:GLN:HG2	1:A:133:ARG:NH1	2.33	0.43
3:C:37:GLY:HA3	3:C:39:TYR:CE2	2.54	0.43
3:C:38:ASN:HB3	3:G:38:ASN:O	2.19	0.43
5:I:19:DA:H2"	5:I:20:DT:H5'	2.01	0.43
5:I:118:DT:H1'	5:I:119:DT:H5'	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:130:DT:H2"	5:I:131:DG:N7	2.34	0.43
5:I:146:DT:H6	5:I:146:DT:H5'	1.84	0.43
1:A:112:HIS:HE1	1:E:121:LYS:HG3	1.84	0.42
3:G:16:THR:O	3:G:19:SER:OG	2.30	0.42
5:I:63:DG:H1'	5:I:64:DT:H5'	2.01	0.42
5:I:98:DG:N2	5:J:195:DC:N3	2.61	0.42
4:D:34:LYS:HB2	5:J:270:DA:OP1	2.19	0.42
5:J:197:DA:C2'	5:J:198:DT:H5"	2.50	0.42
4:D:120:LYS:HB2	4:D:120:LYS:HE2	1.87	0.42
1:A:81:LEU:HD23	1:A:81:LEU:HA	1.72	0.42
1:A:107:ASN:ND2	2:B:42:GLY:O	2.53	0.42
5:I:48:DT:H2"	5:I:49:DC:H5'	2.01	0.42
5:I:38:DT:H2"	5:I:39:DG:C8	2.54	0.42
3:G:54:VAL:HG13	4:H:110:ALA:HB1	2.01	0.42
5:J:181:DA:H2"	5:J:182:DT:H5"	2.02	0.42
5:J:214:DG:H2"	5:J:215:DC:C6	2.54	0.42
5:I:83:DA:H1'	5:I:84:DC:H5'	2.01	0.42
5:I:102:DA:H2"	5:I:103:DG:C8	2.53	0.42
3:C:32:ARG:NH2	4:D:35:GLU:OE1	2.34	0.41
5:I:1:DA:N6	5:J:292:DT:H3	2.18	0.41
5:J:274:DT:C2'	5:J:275:DC:H5"	2.48	0.41
4:D:68:ASP:OD2	2:F:98:TYR:OH	2.28	0.41
5:J:242:DT:H6	5:J:242:DT:H2'	1.75	0.41
2:B:62:LEU:HD23	2:B:62:LEU:HA	1.89	0.41
3:C:39:TYR:HB3	4:D:78:SER:HB2	2.02	0.41
4:D:62:MET:O	4:D:66:VAL:HG23	2.20	0.41
1:A:81:LEU:HD11	2:B:70:VAL:HG13	2.02	0.41
1:E:77:PHE:CZ	2:F:67:ARG:HB2	2.56	0.41
2:F:92:ARG:HB3	2:F:92:ARG:NH1	2.34	0.41
5:I:15:DG:H1'	5:I:16:DC:H5'	2.01	0.41
5:J:198:DT:C2'	5:J:199:DC:H5'	2.49	0.41
3:G:26:PRO:HD3	4:H:40:TYR:CD1	2.56	0.41
2:F:83:ALA:O	2:F:87:VAL:HG23	2.21	0.41
4:H:40:TYR:OH	5:I:121:DG:H5"	2.21	0.41
4:H:102:LEU:HA	4:H:103:PRO:HD3	1.88	0.41
5:J:248:DA:H2"	5:J:249:DG:C8	2.54	0.41
5:I:139:DA:H2"	5:I:140:DT:H5'	2.03	0.41
5:J:206:DC:H2"	5:J:207:DA:C8	2.56	0.41
1:A:70:VAL:HG13	2:B:66:ILE:HD11	2.03	0.40
5:I:37:DT:H2"	5:I:38:DT:H5'	2.02	0.40
5:J:277:DG:H5"	5:J:277:DG:C8	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:30:THR:HB	2:F:32:PRO:HD2	2.03	0.40
4:D:46:LYS:HD3	4:D:46:LYS:HA	1.77	0.40
4:H:88:THR:HG22	5:J:186:DG:OP1	2.20	0.40
1:E:78:ASN:HB3	1:E:81:LEU:HG	2.04	0.40
5:I:57:DA:H2"	5:I:58:DG:C8	2.57	0.40
5:J:249:DG:H1'	5:J:250:DT:H5"	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:79:THR:OG1	5:J:285:DA:OP1[3_647]	2.17	0.03

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	95/138 (69%)	95 (100%)	0	0	100	100
1	E	94/138 (68%)	94 (100%)	0	0	100	100
2	B	81/106 (76%)	80 (99%)	1 (1%)	0	100	100
2	F	77/106 (73%)	76 (99%)	1 (1%)	0	100	100
3	C	103/133 (77%)	102 (99%)	1 (1%)	0	100	100
3	G	102/133 (77%)	101 (99%)	1 (1%)	0	100	100
4	D	90/129 (70%)	89 (99%)	1 (1%)	0	100	100
4	H	90/129 (70%)	89 (99%)	1 (1%)	0	100	100
All	All	732/1012 (72%)	726 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	83/113 (74%)	83 (100%)	0	100	100
1	E	82/113 (73%)	82 (100%)	0	100	100
2	B	68/81 (84%)	68 (100%)	0	100	100
2	F	64/81 (79%)	64 (100%)	0	100	100
3	C	83/102 (81%)	82 (99%)	1 (1%)	71	92
3	G	83/102 (81%)	82 (99%)	1 (1%)	71	92
4	D	78/107 (73%)	78 (100%)	0	100	100
4	H	78/107 (73%)	78 (100%)	0	100	100
All	All	619/806 (77%)	617 (100%)	2 (0%)	92	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	35	ARG
3	G	118	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides 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	97/138 (70%)	0.11	1 (1%)	82	77	37, 55, 85, 132	0
1	E	96/138 (69%)	0.17	2 (2%)	63	54	47, 75, 109, 122	0
2	B	83/106 (78%)	0.17	0	100	100	36, 50, 70, 85	0
2	F	79/106 (74%)	0.16	2 (2%)	57	47	46, 70, 97, 128	0
3	C	105/133 (78%)	0.07	1 (0%)	82	77	50, 69, 105, 119	0
3	G	104/133 (78%)	0.10	1 (0%)	82	77	38, 56, 85, 116	0
4	D	92/129 (71%)	0.19	1 (1%)	80	75	46, 68, 93, 127	0
4	H	92/129 (71%)	0.13	0	100	100	38, 61, 83, 108	0
5	I	146/146 (100%)	-0.11	2 (1%)	75	70	77, 131, 169, 212	0
5	J	146/146 (100%)	-0.10	2 (1%)	75	70	71, 131, 166, 200	0
All	All	1040/1304 (79%)	0.07	12 (1%)	79	73	36, 71, 155, 212	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	J	147	DA	7.2
5	I	146	DT	6.0
1	E	38	HIS	3.9
2	F	24	ASP	2.8
1	E	40	TYR	2.5
3	C	15	LYS	2.5
5	J	148	DT	2.4
3	G	118	LYS	2.4
5	I	1	DA	2.4
2	F	102	GLY	2.2
1	A	80	ASP	2.1
4	D	124	ALA	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 monosaccharides 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.