



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 05:03 PM UTC

PDB ID : 5GSE / pdb_00005gse
Title : Crystal structure of unusual nucleosome
Authors : Kato, D.; Osakabe, A.; Arimura, Y.; Park, S.Y.; Kurumizaka, H.
Deposited on : 2016-08-16
Resolution : 3.14 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

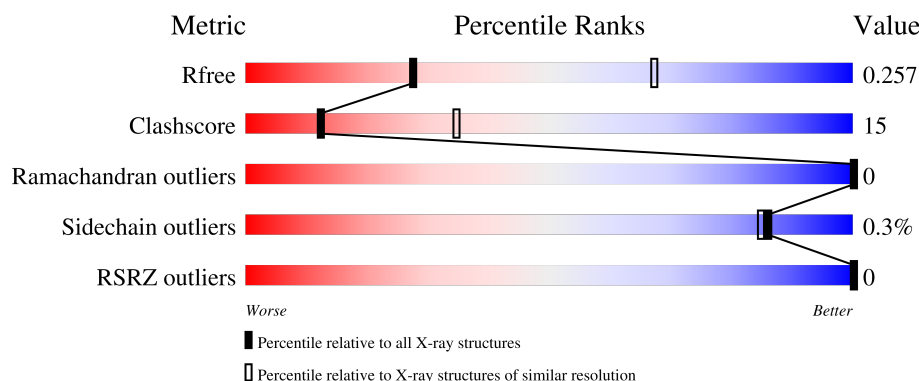
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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.



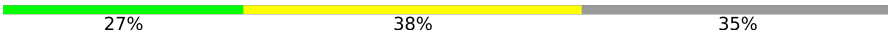
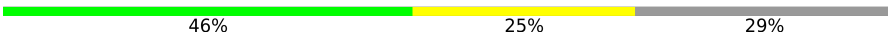
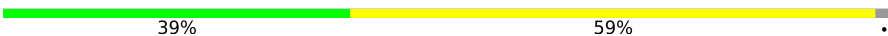
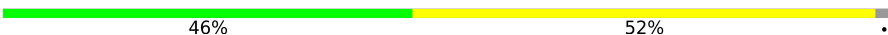
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	139	
1	E	139	
1	K	139	
1	O	139	
2	B	106	

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Mol	Chain	Length	Quality of chain
2	F	106	
2	L	106	
2	P	106	
3	C	133	
3	G	133	
3	M	133	
4	D	129	
4	H	129	
4	N	129	
5	I	250	
6	J	250	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 20244 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	96	Total	C	N	O	S	0	0	0
			794	500	154	136	4			
1	E	97	Total	C	N	O	S	0	0	0
			801	505	155	137	4			
1	K	88	Total	C	N	O	S	0	0	0
			725	457	138	126	4			
1	O	96	Total	C	N	O	S	0	0	0
			794	500	154	136	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P68431
A	-2	SER	-	expression tag	UNP P68431
A	-1	HIS	-	expression tag	UNP P68431
E	-3	GLY	-	expression tag	UNP P68431
E	-2	SER	-	expression tag	UNP P68431
E	-1	HIS	-	expression tag	UNP P68431
K	-3	GLY	-	expression tag	UNP P68431
K	-2	SER	-	expression tag	UNP P68431
K	-1	HIS	-	expression tag	UNP P68431
O	-3	GLY	-	expression tag	UNP P68431
O	-2	SER	-	expression tag	UNP P68431
O	-1	HIS	-	expression tag	UNP P68431

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	82	Total	C	N	O	S	0	0	0
			653	412	127	113	1			
2	F	79	Total	C	N	O	S	0	0	0
			633	399	124	109	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	69	Total	C	N	O	S	0	0	0
			553	349	108	95	1			
2	P	81	Total	C	N	O	S	0	0	0
			648	410	126	111	1			

There are 12 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
L	-3	GLY	-	expression tag	UNP P62805
L	-2	SER	-	expression tag	UNP P62805
L	-1	HIS	-	expression tag	UNP P62805
P	-3	GLY	-	expression tag	UNP P62805
P	-2	SER	-	expression tag	UNP P62805
P	-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
3	C	102	Total	C	N	O		0	0	0
			787	496	153	138				
3	G	109	Total	C	N	O		0	0	0
			840	529	166	145				
3	M	106	Total	C	N	O		0	0	0
			815	514	159	142				

There are 9 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
M	-3	GLY	-	expression tag	UNP P04908
M	-2	SER	-	expression tag	UNP P04908

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Chain	Residue	Modelled	Actual	Comment	Reference
M	-1	HIS	-	expression tag	UNP P04908

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	92	Total	C	N	O	S	0	0	0
			720	453	129	136	2			
4	H	91	Total	C	N	O	S	0	0	0
			708	447	125	134	2			
4	N	93	Total	C	N	O	S	0	0	0
			725	456	130	137	2			

There are 9 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
N	-3	GLY	-	expression tag	UNP P06899
N	-2	SER	-	expression tag	UNP P06899
N	-1	HIS	-	expression tag	UNP P06899

- Molecule 5 is a DNA chain called DNA (250-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	245	Total	C	N	O	P	0	0	0
			5055	2393	948	1469	245			

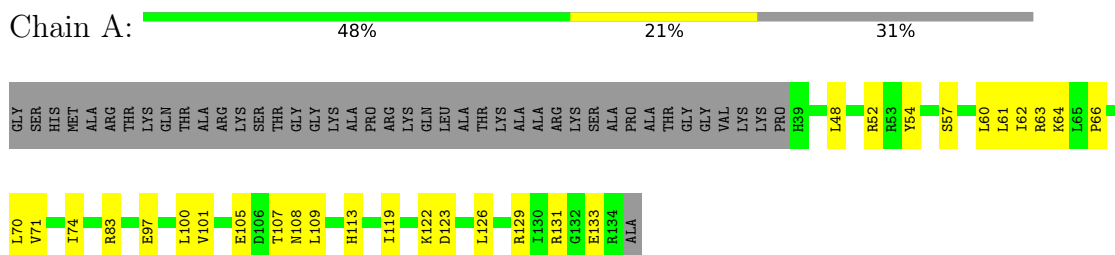
- Molecule 6 is a DNA chain called DNA (250-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	J	245	Total	C	N	O	P	0	0	0
			4993	2372	905	1471	245			

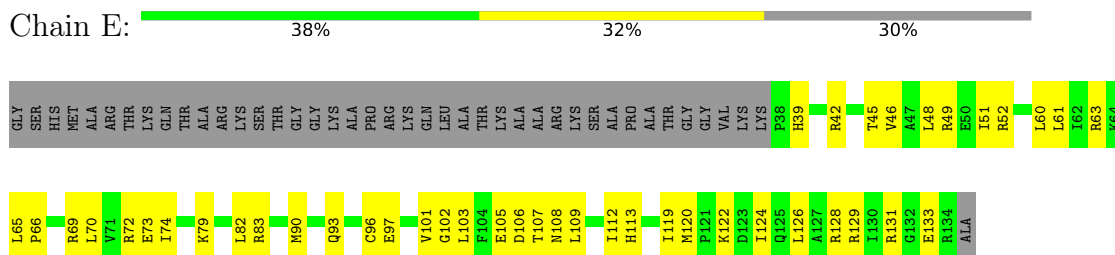
3 Residue-property plots

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.

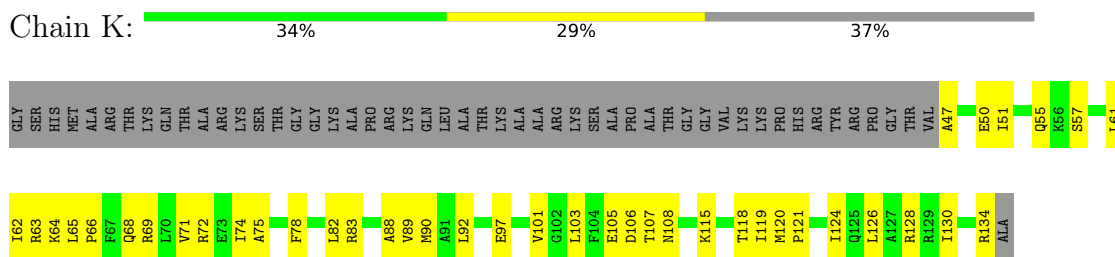
• Molecule 1: Histone H3.1



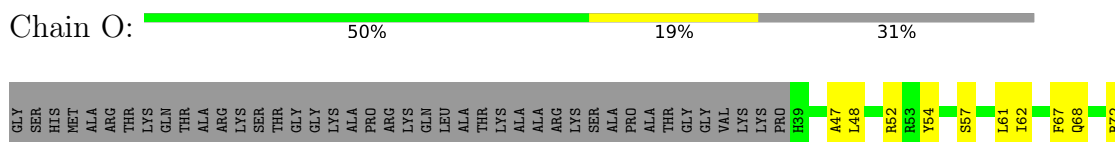
• Molecule 1: Histone H3.1



• Molecule 1: Histone H3.1

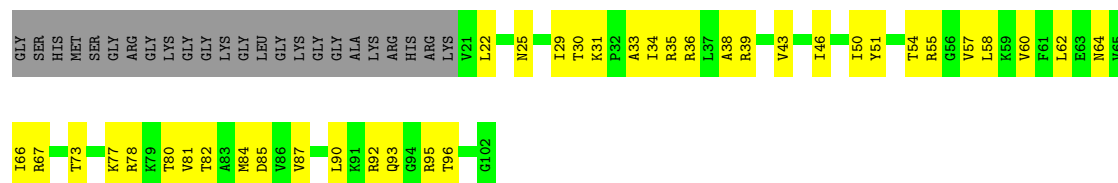


• Molecule 1: Histone H3.1





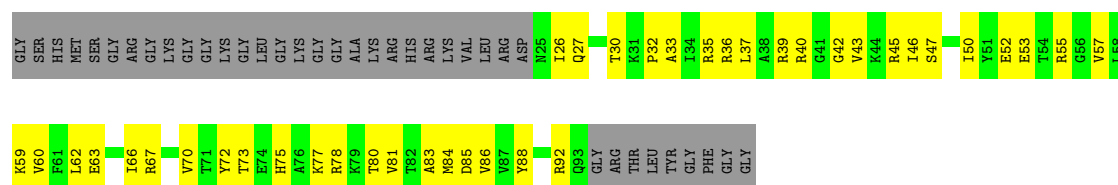
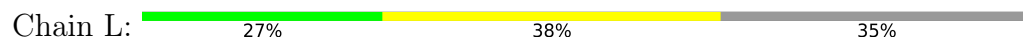
• Molecule 2: Histone H4



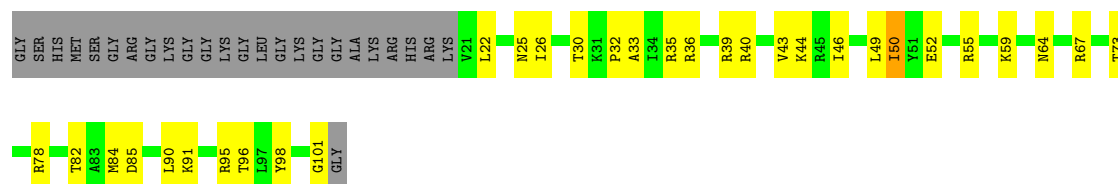
• Molecule 2: Histone H4



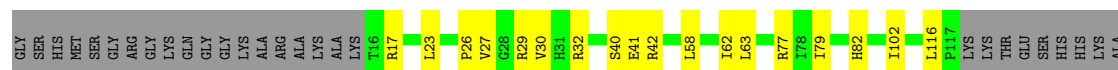
• Molecule 2: Histone H4



• Molecule 2: Histone H4



• Molecule 3: Histone H2A type 1-B/E



LYS
GLY
LYS

• Molecule 3: Histone H2A type 1-B/E

Chain G:  57% 25% 18%

GLY SER HIS MET SER GLY ARG GLY LYS GLN GLY K9
T16 R17 S18 S19 R20 Q24 F25 P26 V27 G28 R29 V30 H31 G37 N38 Y39 R42 P48 V49 V54 L55 E56 V57 L58 T59 I62 L63 M68 R71 R77 I78 I79 P80 R81 H82 L83 Q84
R88 V100 T101 I102 G106 P117 LYS LYS THR GLU SER HIS HIS LYS LYS ALA GLY LYS

• Molecule 3: Histone H2A type 1-B/E

Chain M:  62% 17% 20%

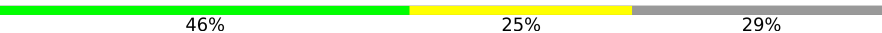
GLY SER HIS MET SER GLY ARG GLY LYS GLN GLY LYS ALA A12 T16 R17 P26 V27 G28 R29 V30 Y39 S40 E41 R42 G43 G44 L51 V54 L58 L63 E64 L65 A66 G67 N68 N73 K74 K75 I79 P80 L83 L83 P117 LYS
THR GLU SER HIS LYS ALA LYS LYS

• Molecule 4: Histone H2B type 1-J

Chain D:  53% 18% 29%

GLY SER HIS MET PRO GLU PRO ALA LYS SER SER PRO ALA ALA PRO PRO LYS LYS GLY SER LYS LYS ALA VAL THR LYS LYS ALA GLN LYS LYS ASP LYS LYS ARG ARG S32 E35 Y40 V41 V44 L45 I61 M62 D68 R72 N84 K85 R86 S87 T88 I89 T90 S91
R92 E93 I94 Q95 T96 A97 V98 L101 L102 K108 V111 Y121 T122 S123 ALA LYS

• Molecule 4: Histone H2B type 1-J

Chain H:  46% 25% 29%

GLY SER HIS MET PRO GLU PRO ALA LYS SER SER PRO ALA ALA PRO PRO LYS LYS GLY SER LYS LYS ALA VAL THR LYS LYS ALA GLN LYS LYS ASP LYS LYS ARG ARG S32 K34 E35 S36 Y40 K43 Y44 L45 I54 S55 A58 T61 M62 N63 V66 I69
I73 S78 T88 I89 T90 S91 E93 E94 Q95 T96 R99 L100 L101 L102 E105 K108 H109 A110 V111 S112 E113 G114 V118 A124 LYS

• Molecule 4: Histone H2B type 1-J

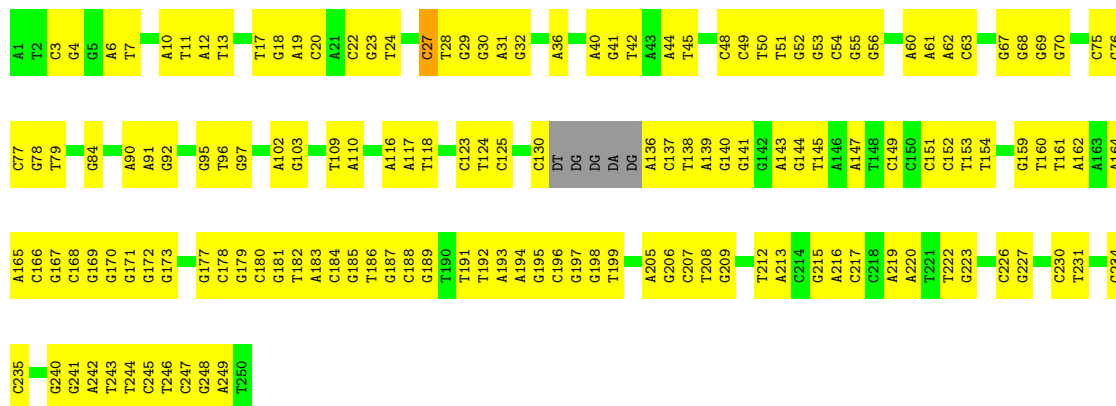
Chain N:  50% 22% 28%

GLY SER HIS MET PRO GLU PRO ALA LYS SER SER PRO ALA ALA PRO PRO LYS LYS GLY SER LYS LYS ALA VAL THR LYS LYS ALA GLN LYS LYS ASP LYS LYS ARG ARG S32 R33 K34 E35 Y40 V44 L45 V48 G53 S56 F58 D68 S78 Y83 R86



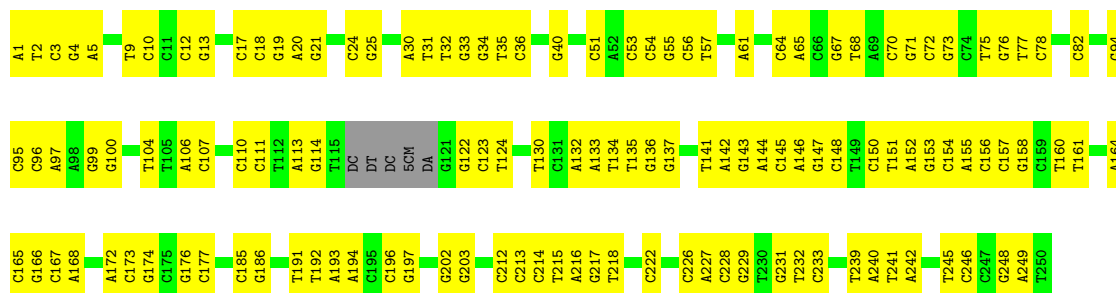
• Molecule 5: DNA (250-MER)

Chain I: 39% 59%



• Molecule 6: DNA (250-MER)

Chain J: 46% 52%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	90.58Å 101.80Å 102.43Å 119.30° 106.51° 91.36°	Depositor
Resolution (Å)	49.66 – 3.14 49.66 – 3.14	Depositor EDS
% Data completeness (in resolution range)	91.0 (49.66-3.14) 90.9 (49.66-3.14)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.197 , 0.255 0.199 , 0.257	Depositor DCC
R_{free} test set	2369 reflections (4.46%)	wwPDB-VP
Wilson B-factor (Å ²)	92.8	Xtriage
Anisotropy	0.376	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.024 for -h,k,-k-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20244	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 5CM

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	0.22	0/805	0.52	0/1079
1	E	0.25	0/813	0.58	0/1090
1	K	0.28	0/733	0.67	0/981
1	O	0.25	0/805	0.53	0/1079
2	B	0.27	0/660	0.58	0/883
2	F	0.22	0/640	0.49	0/857
2	L	0.19	0/558	0.50	0/748
2	P	0.27	0/655	0.57	0/878
3	C	0.24	0/797	0.51	0/1078
3	G	0.22	0/850	0.52	0/1146
3	M	0.23	0/825	0.53	0/1114
4	D	0.19	0/731	0.44	0/983
4	H	0.22	0/719	0.53	0/968
4	N	0.23	0/736	0.51	0/990
5	I	0.24	0/5631	0.55	0/8693
6	J	0.24	0/5570	0.51	0/8582
All	All	0.24	0/21528	0.53	0/31149

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

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	794	0	831	25	0
1	E	801	0	839	45	0
1	K	725	0	763	42	0
1	O	794	0	831	25	0
2	B	653	0	696	41	0
2	F	633	0	673	30	0
2	L	553	0	598	41	0
2	P	648	0	693	33	0
3	C	787	0	835	19	0
3	G	840	0	902	30	0
3	M	815	0	871	20	0
4	D	720	0	740	20	0
4	H	708	0	727	28	0
4	N	725	0	745	25	0
5	I	5055	0	2754	146	0
6	J	4993	0	2753	120	0
All	All	20244	0	16251	557	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 15.

The worst 5 of 557 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:K:130:ILE:HD11	1:O:130:ILE:HB	1.54	0.90
5:I:84:DG:N2	6:J:167:DC:O2	2.04	0.89
6:J:1:DA:H2''	6:J:2:DT:H71	1.60	0.82
3:M:79:ILE:HG13	3:M:80:PRO:HD2	1.62	0.80
5:I:185:DG:OP2	2:L:35:ARG:NH1	2.15	0.78

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	94/139 (68%)	93 (99%)	1 (1%)	0	100	100
1	E	95/139 (68%)	90 (95%)	5 (5%)	0	100	100
1	K	86/139 (62%)	82 (95%)	4 (5%)	0	100	100
1	O	94/139 (68%)	91 (97%)	3 (3%)	0	100	100
2	B	80/106 (76%)	76 (95%)	4 (5%)	0	100	100
2	F	77/106 (73%)	74 (96%)	3 (4%)	0	100	100
2	L	67/106 (63%)	65 (97%)	2 (3%)	0	100	100
2	P	79/106 (74%)	77 (98%)	2 (2%)	0	100	100
3	C	100/133 (75%)	96 (96%)	4 (4%)	0	100	100
3	G	107/133 (80%)	104 (97%)	3 (3%)	0	100	100
3	M	104/133 (78%)	101 (97%)	3 (3%)	0	100	100
4	D	90/129 (70%)	88 (98%)	2 (2%)	0	100	100
4	H	89/129 (69%)	87 (98%)	2 (2%)	0	100	100
4	N	91/129 (70%)	89 (98%)	2 (2%)	0	100	100
All	All	1253/1766 (71%)	1213 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	84/113 (74%)	84 (100%)	0	100	100
1	E	85/113 (75%)	85 (100%)	0	100	100
1	K	77/113 (68%)	76 (99%)	1 (1%)	61	73
1	O	84/113 (74%)	84 (100%)	0	100	100
2	B	67/81 (83%)	67 (100%)	0	100	100
2	F	65/81 (80%)	64 (98%)	1 (2%)	57	71
2	L	58/81 (72%)	58 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	P	67/81 (83%)	66 (98%)	1 (2%)	57	71
3	C	81/102 (79%)	81 (100%)	0	100	100
3	G	85/102 (83%)	85 (100%)	0	100	100
3	M	83/102 (81%)	83 (100%)	0	100	100
4	D	79/107 (74%)	79 (100%)	0	100	100
4	H	77/107 (72%)	77 (100%)	0	100	100
4	N	79/107 (74%)	79 (100%)	0	100	100
All	All	1071/1403 (76%)	1068 (100%)	3 (0%)	86	85

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	73	THR
1	K	115	LYS
2	P	50	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
4	N	84	ASN
1	O	125	GLN
3	G	104	GLN
4	H	95	GLN
1	K	113	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

3 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	5CM	I	130	6,5	18,21,22	1.66	2 (11%)	24,30,33	1.23	2 (8%)
6	5CM	J	222	6,5	18,21,22	1.75	3 (16%)	24,30,33	1.27	2 (8%)
5	5CM	I	27	6,5	18,21,22	1.81	3 (16%)	24,30,33	1.37	5 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	5CM	I	130	6,5	-	0/7/21/22	0/2/2/2
6	5CM	J	222	6,5	-	4/7/21/22	0/2/2/2
5	5CM	I	27	6,5	-	4/7/21/22	0/2/2/2

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	27	5CM	C5-C4	6.36	1.48	1.44
6	J	222	5CM	C5-C4	6.13	1.48	1.44
5	I	130	5CM	C5-C4	5.74	1.48	1.44
5	I	27	5CM	C6-C5	3.07	1.39	1.34
5	I	130	5CM	C6-C5	2.93	1.39	1.34

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	222	5CM	C5-C6-N1	-4.06	118.91	123.31
5	I	130	5CM	C5-C6-N1	-3.92	119.05	123.31
5	I	27	5CM	C5-C6-N1	-3.67	119.32	123.31
5	I	27	5CM	C5-C4-N3	-2.65	119.03	121.75
5	I	27	5CM	O4'-C1'-N1	2.33	112.00	107.86

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	I	27	5CM	C2'-C1'-N1-C6

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Mol	Chain	Res	Type	Atoms
5	I	27	5CM	O4'-C1'-N1-C6
6	J	222	5CM	C2'-C1'-N1-C6
5	I	27	5CM	C2'-C1'-N1-C2
6	J	222	5CM	O4'-C1'-N1-C6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	27	5CM	1	0

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	96/139 (69%)	-0.63	0 100 100	48, 64, 84, 103	0
1	E	97/139 (69%)	-0.48	0 100 100	59, 74, 100, 115	0
1	K	88/139 (63%)	-0.47	0 100 100	61, 75, 103, 129	0
1	O	96/139 (69%)	-0.43	0 100 100	58, 75, 100, 121	0
2	B	82/106 (77%)	-0.42	0 100 100	51, 64, 84, 92	0
2	F	79/106 (74%)	-0.45	0 100 100	53, 73, 90, 111	0
2	L	69/106 (65%)	-0.45	0 100 100	55, 76, 90, 95	0
2	P	81/106 (76%)	-0.41	0 100 100	53, 73, 101, 109	0
3	C	102/133 (76%)	-0.49	0 100 100	61, 73, 94, 103	0
3	G	109/133 (81%)	-0.57	0 100 100	57, 76, 97, 118	0
3	M	106/133 (79%)	-0.55	0 100 100	55, 71, 94, 115	0
4	D	92/129 (71%)	-0.47	0 100 100	57, 75, 90, 106	0
4	H	91/129 (70%)	-0.43	0 100 100	63, 79, 94, 109	0
4	N	93/129 (72%)	-0.61	0 100 100	51, 72, 91, 118	0
5	I	243/250 (97%)	-0.53	0 100 100	78, 115, 145, 162	0
6	J	244/250 (97%)	-0.53	0 100 100	74, 117, 145, 159	0
All	All	1768/2266 (78%)	-0.50	0 100 100	48, 79, 133, 162	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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($\text{\AA}^2$)	Q<0.9
5	5CM	I	130	20/21	0.63	0.11	98,124,138,154	0
6	5CM	J	222	20/21	0.82	0.10	99,113,121,123	0
5	5CM	I	27	20/21	0.91	0.08	93,111,123,132	0

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.