



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 01:08 am BST

PDB ID : 5GSE
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 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

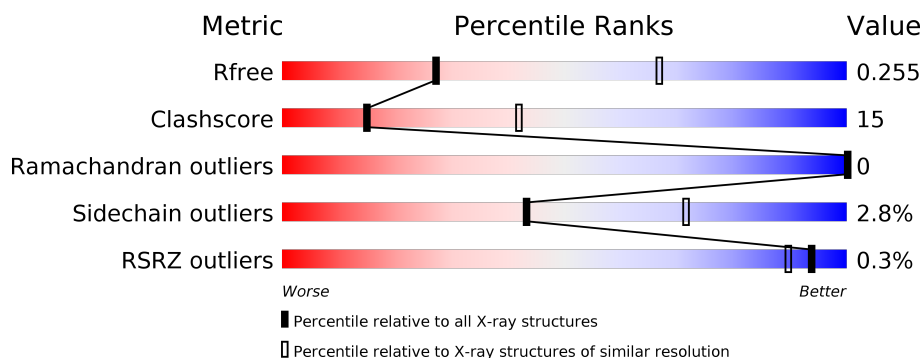
1 Overall quality at a glance

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.









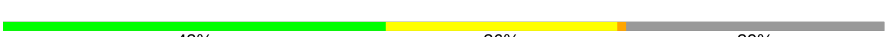



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1626 (3.18-3.10)
Clashscore	141614	1735 (3.18-3.10)
Ramachandran outliers	138981	1677 (3.18-3.10)
Sidechain outliers	138945	1677 (3.18-3.10)
RSRZ outliers	127900	1588 (3.18-3.10)

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 on 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	<div> <div style="width: 48%; background-color: green;"></div> <div style="width: 20%; background-color: yellow;"></div> <div style="width: 31%; background-color: grey;"></div> </div> <div>48% 20% • 31%</div>
1	E	139	<div> <div style="width: 39%; background-color: green;"></div> <div style="width: 29%; background-color: yellow;"></div> <div style="width: 30%; background-color: grey;"></div> </div> <div>39% 29% •• 30%</div>
1	K	139	<div> <div style="width: 36%; background-color: green;"></div> <div style="width: 25%; background-color: yellow;"></div> <div style="width: 37%; background-color: grey;"></div> </div> <div>36% 25% • 37%</div>
1	O	139	<div> <div style="width: 50%; background-color: green;"></div> <div style="width: 18%; background-color: yellow;"></div> <div style="width: 31%; background-color: grey;"></div> </div> <div>% 50% 18% • 31%</div>
2	B	106	<div> <div style="width: 42%; background-color: green;"></div> <div style="width: 33%; background-color: yellow;"></div> <div style="width: 23%; background-color: grey;"></div> </div> <div>42% 33% • 23%</div>
2	F	106	<div> <div style="width: 45%; background-color: green;"></div> <div style="width: 28%; background-color: yellow;"></div> <div style="width: 25%; background-color: grey;"></div> </div> <div>% 45% 28% • 25%</div>

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Mol	Chain	Length	Quality of chain
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			

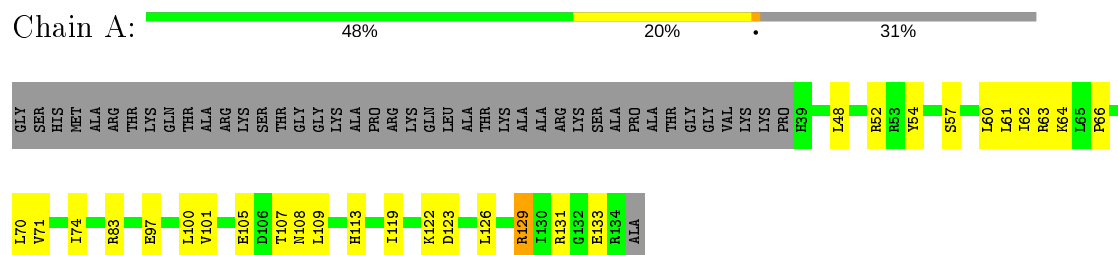
- Molecule 6 is a DNA chain called DNA (250-MER).

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

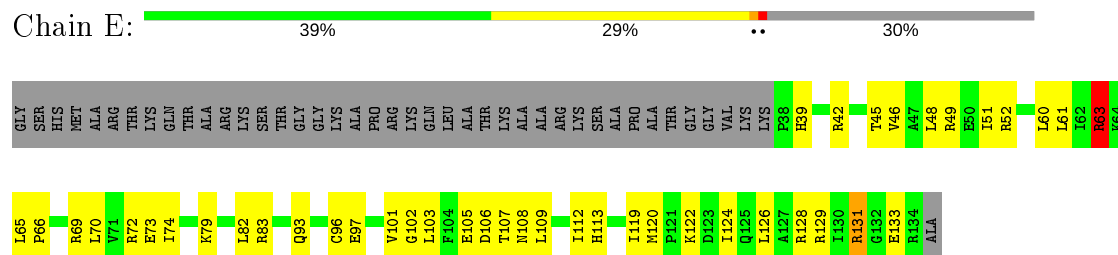
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

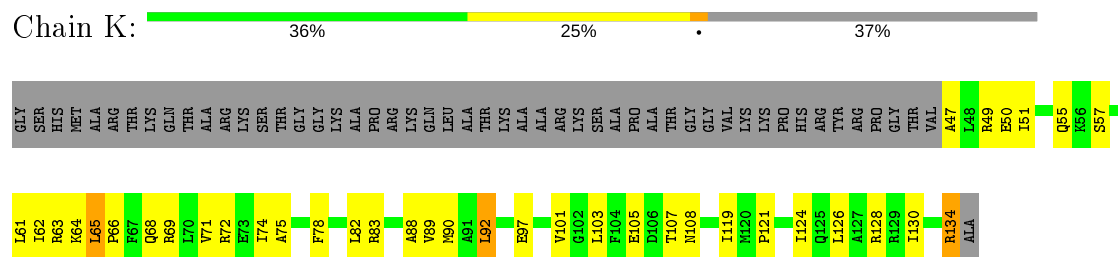
• Molecule 1: Histone H3.1



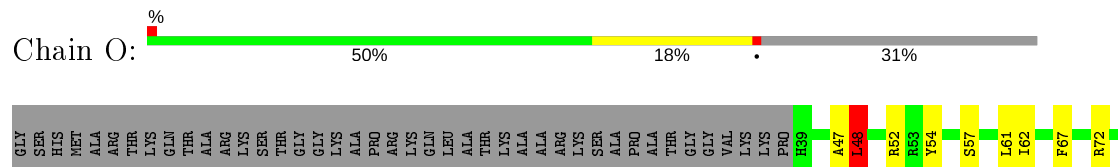
• Molecule 1: Histone H3.1



• Molecule 1: Histone H3.1

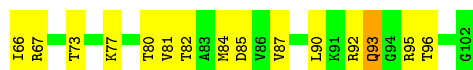
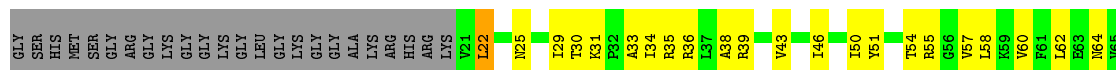


• Molecule 1: Histone H3.1

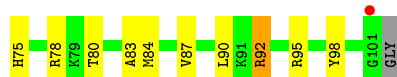




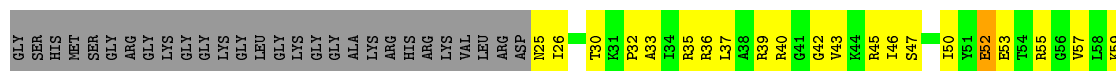
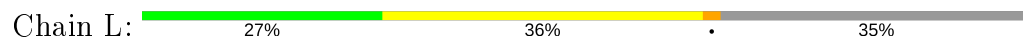
• Molecule 2: Histone H4



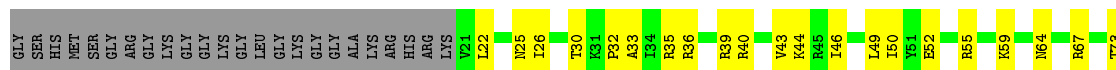
• Molecule 2: Histone H4



• Molecule 2: Histone H4

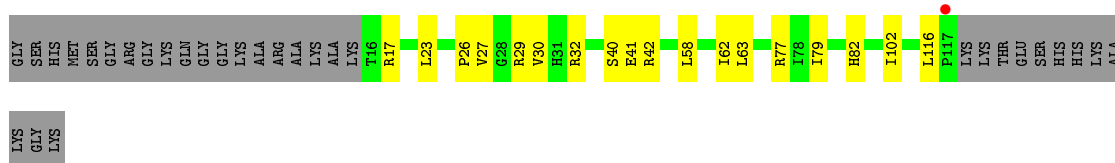


• Molecule 2: Histone H4

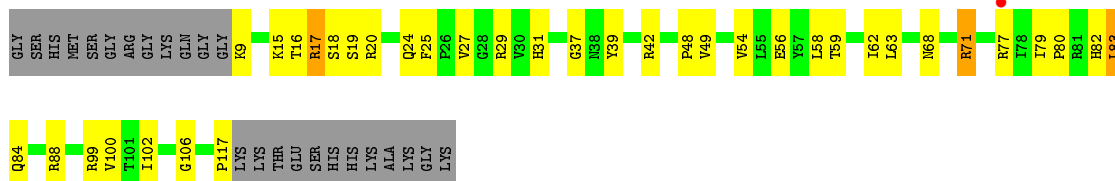


• Molecule 3: Histone H2A type 1-B/E

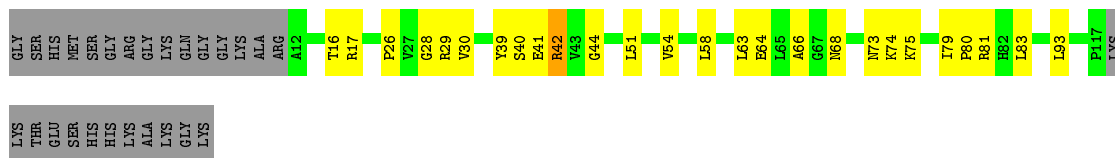




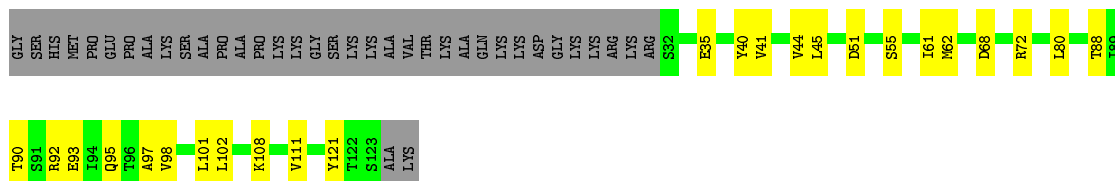
• Molecule 3: Histone H2A type 1-B/E



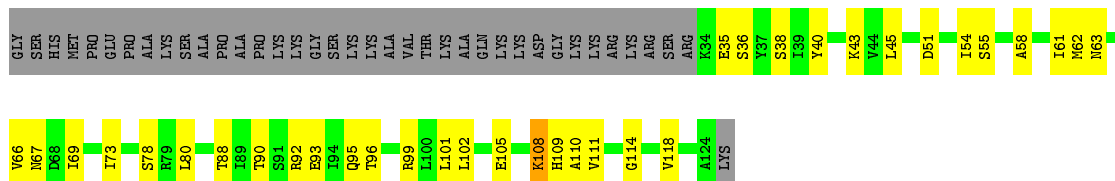
• Molecule 3: Histone H2A type 1-B/E



• Molecule 4: Histone H2B type 1-J

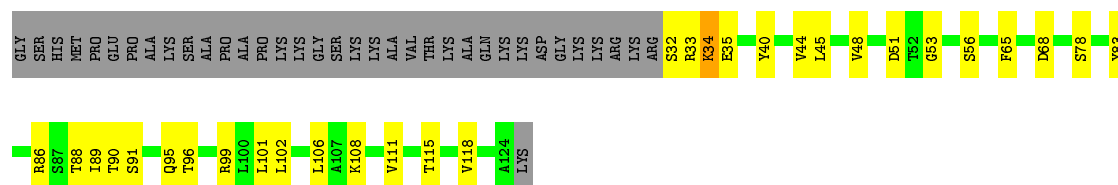


• Molecule 4: Histone H2B type 1-J



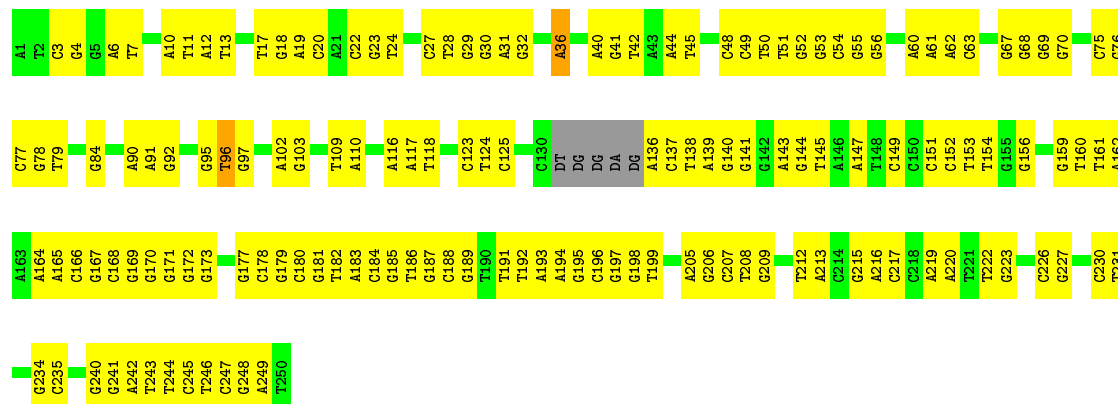
• Molecule 4: Histone H2B type 1-J





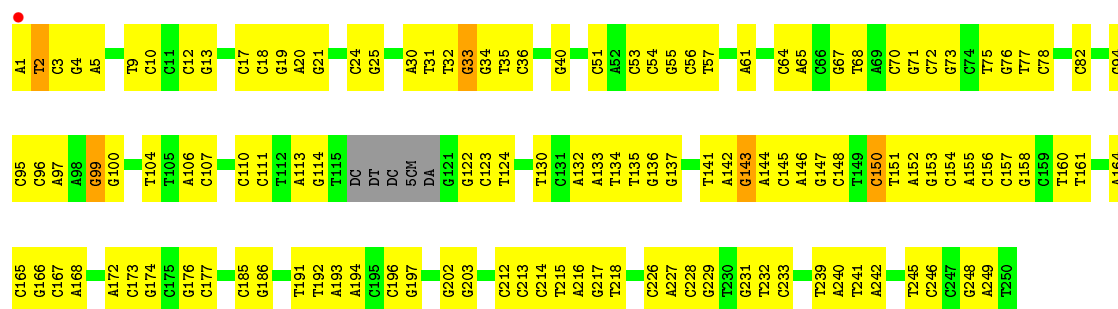
• Molecule 5: DNA (250-MER)

Chain I: ..



• Molecule 6: DNA (250-MER)

Chain J: ..



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) 91.0 (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 0.197 , 0.255	Depositor DCC
R_{free} test set	2359 reflections (4.95%)	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)	0.29 , 56.7	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 [i](#)

5.1 Standard geometry [i](#)

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	0.29	0/805	0.53	0/1079
1	E	0.31	0/813	0.63	1/1090 (0.1%)
1	K	0.33	0/733	0.70	2/981 (0.2%)
1	O	0.31	0/805	0.55	1/1079 (0.1%)
2	B	0.34	0/660	0.56	0/883
2	F	0.29	0/640	0.53	0/857
2	L	0.26	0/558	0.47	0/748
2	P	0.31	0/655	0.60	0/878
3	C	0.29	0/797	0.54	0/1078
3	G	0.29	0/850	0.59	1/1146 (0.1%)
3	M	0.31	0/825	0.56	1/1114 (0.1%)
4	D	0.29	0/731	0.49	1/983 (0.1%)
4	H	0.29	0/719	0.52	0/968
4	N	0.32	0/736	0.53	0/990
5	I	0.62	0/5631	1.00	3/8693 (0.0%)
6	J	0.60	0/5570	1.01	6/8582 (0.1%)
All	All	0.48	0/21528	0.84	16/31149 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	92	LEU	CA-CB-CG	7.23	131.94	115.30
5	I	96	DT	O4'-C4'-C3'	-7.12	101.65	104.50
1	E	63	ARG	NE-CZ-NH2	-6.99	116.81	120.30
1	K	65	LEU	CA-CB-CG	6.38	129.97	115.30
3	G	83	LEU	CA-CB-CG	6.31	129.82	115.30
5	I	36	DA	O4'-C4'-C3'	-5.88	102.15	104.50
1	O	48	LEU	CA-CB-CG	5.71	128.43	115.30
4	D	80	LEU	CA-CB-CG	5.60	128.18	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	150	DC	O4'-C1'-N1	5.48	111.84	108.00
6	J	2	DT	N3-C4-O4	5.45	123.17	119.90
6	J	143	DG	O4'-C1'-N9	5.34	111.73	108.00
5	I	156	DG	O4'-C4'-C3'	-5.30	102.38	104.50
6	J	2	DT	C5-C4-O4	-5.28	121.21	124.90
6	J	99	DG	O4'-C1'-N9	5.19	111.63	108.00
6	J	33	DG	O4'-C4'-C3'	-5.02	102.49	104.50
3	M	42	ARG	NE-CZ-NH1	-5.01	117.79	120.30

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	44	0
1	K	725	0	763	40	0
1	O	794	0	831	22	0
2	B	653	0	696	39	0
2	F	633	0	673	30	0
2	L	553	0	598	40	0
2	P	648	0	693	31	0
3	C	787	0	835	19	0
3	G	840	0	902	29	0
3	M	815	0	871	21	0
4	D	720	0	740	19	0
4	H	708	0	727	27	0
4	N	725	0	745	25	0
5	I	5055	0	2754	145	0
6	J	4993	0	2753	121	0
All	All	20244	0	16251	546	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.

All (546) 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
3:M:79:ILE:HG13	3:M:80:PRO:HD2	1.62	0.80
1:K:51:ILE:O	1:K:55:GLN:HB2	1.83	0.78
5:I:185:DG:OP2	2:L:35:ARG:NH1	2.15	0.78
3:C:42:ARG:HB2	4:D:88:THR:HG22	1.66	0.78
6:J:106:DA:H2''	6:J:107:DC:H5''	1.67	0.76
1:E:83:ARG:HB2	2:F:80:THR:HG22	1.66	0.76
1:E:42:ARG:NH2	6:J:172:DA:OP1	2.14	0.75
1:O:61:LEU:O	2:P:36:ARG:NH1	2.21	0.74
5:I:164:DA:OP1	2:P:36:ARG:NH2	2.21	0.73
5:I:144:DG:H1	6:J:107:DC:H5	1.38	0.72
1:E:79:LYS:HB3	1:E:82:LEU:HD21	1.72	0.72
1:O:131:ARG:NH1	1:O:133:GLU:OE2	2.22	0.72
3:C:79:ILE:H	3:C:82:HIS:HD1	1.38	0.71
3:M:42:ARG:HB2	4:N:88:THR:HG22	1.73	0.70
6:J:245:DT:H2'	6:J:246:DC:C6	2.27	0.70
5:I:194:DA:OP2	1:K:69:ARG:NH2	2.25	0.70
6:J:77:DT:H1'	6:J:78:DC:H5'	1.73	0.69
3:M:93:LEU:HD12	4:N:106:LEU:HD21	1.73	0.69
4:H:90:THR:HG23	4:H:92:ARG:H	1.58	0.69
1:O:121:PRO:HA	1:O:124:ILE:HG22	1.75	0.69
2:P:30:THR:OG1	2:P:32:PRO:HD2	1.93	0.68
4:H:73:ILE:HD13	4:H:101:LEU:HD12	1.75	0.68
2:B:29:ILE:O	2:B:55:ARG:NH1	2.27	0.68
5:I:152:DC:H2''	5:I:153:DT:H71	1.75	0.68
6:J:31:DT:H2'	6:J:32:DT:H71	1.76	0.68
6:J:24:DC:H2''	6:J:25:DG:C8	2.28	0.67
6:J:64:DC:H2''	6:J:65:DA:H5'	1.76	0.67
6:J:61:DA:OP1	2:L:36:ARG:NH2	2.27	0.67
2:P:64:ASN:HA	2:P:67:ARG:HG2	1.77	0.67
1:E:83:ARG:HD2	6:J:154:DC:H5'	1.76	0.67
4:H:108:LYS:O	4:H:111:VAL:HG12	1.95	0.66
5:I:181:DG:H2''	5:I:182:DT:H72	1.78	0.65
3:G:59:THR:O	3:G:63:LEU:HB2	1.97	0.64
5:I:159:DG:H2'	5:I:160:DT:C6	2.33	0.64
2:B:50:ILE:O	2:B:54:THR:HG22	1.98	0.64
1:O:119:ILE:HD13	2:P:43:VAL:HG13	1.80	0.64
3:M:40:SER:HB2	4:N:89:ILE:HG13	1.81	0.63
4:D:90:THR:HG23	4:D:92:ARG:H	1.63	0.63
2:P:39:ARG:NH1	2:P:44:LYS:O	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:29:ARG:HH21	4:N:40:TYR:HE2	1.46	0.63
2:B:64:ASN:OD1	2:B:67:ARG:NH2	2.32	0.63
3:C:29:ARG:HH21	4:D:40:TYR:HE2	1.46	0.63
5:I:6:DA:H2"	5:I:7:DT:H5'	1.80	0.62
1:A:101:VAL:O	1:A:105:GLU:HG3	2.00	0.62
2:B:36:ARG:NH2	5:I:61:DA:OP1	2.32	0.62
6:J:185:DC:H2"	6:J:186:DG:C8	2.35	0.62
5:I:159:DG:H2"	5:I:160:DT:O5'	1.99	0.62
1:A:83:ARG:HB2	2:B:80:THR:HG22	1.80	0.62
2:F:26:ILE:O	2:F:29:ILE:HG12	2.00	0.62
3:G:42:ARG:HB2	4:H:88:THR:HG22	1.81	0.62
2:F:29:ILE:O	2:F:55:ARG:NH1	2.33	0.61
5:I:168:DC:H2"	5:I:169:DG:C8	2.35	0.61
2:B:34:ILE:HG21	2:B:54:THR:HG21	1.81	0.61
6:J:202:DG:H2"	6:J:203:DG:N7	2.15	0.61
6:J:51:DC:P	1:K:72:ARG:HH22	2.23	0.61
1:K:108:ASN:HA	1:K:119:ILE:HD11	1.83	0.60
1:K:101:VAL:O	1:K:105:GLU:HG3	2.02	0.60
5:I:160:DT:H4'	5:I:161:DT:OP1	2.02	0.60
5:I:51:DT:H2"	5:I:52:DG:H5"	1.84	0.60
4:N:115:THR:O	4:N:118:VAL:HG12	2.01	0.60
5:I:182:DT:H2"	5:I:183:DA:C8	2.37	0.60
5:I:245:DC:H2"	5:I:246:DT:H71	1.84	0.60
5:I:36:DA:C8	5:I:36:DA:H5'	2.37	0.60
4:H:45:LEU:HD22	4:H:54:ILE:HG13	1.84	0.59
3:M:73:ASN:O	3:M:75:LYS:HG3	2.02	0.59
1:O:62:ILE:HG22	2:P:33:ALA:HB1	1.85	0.59
2:F:30:THR:HG21	6:J:164:DA:H5"	1.83	0.59
2:P:73:THR:OG1	2:P:85:ASP:OD2	2.14	0.59
6:J:64:DC:H2'	6:J:65:DA:C8	2.38	0.59
6:J:31:DT:H2"	6:J:32:DT:H5'	1.85	0.59
2:B:84:MET:HA	2:B:87:VAL:HG12	1.84	0.59
5:I:169:DG:H2"	5:I:170:DG:C8	2.37	0.59
2:P:78:ARG:NH2	2:P:85:ASP:OD2	2.36	0.59
2:F:75:HIS:HB2	4:H:96:THR:HG21	1.84	0.59
4:N:32:SER:O	4:N:34:LYS:NZ	2.35	0.59
1:E:107:THR:HG21	1:E:124:ILE:HG12	1.85	0.59
5:I:247:DC:H2"	5:I:248:DG:C8	2.37	0.59
3:C:23:LEU:HD11	3:C:27:VAL:HG23	1.85	0.58
6:J:20:DA:OP2	4:N:56:SER:OG	2.19	0.58
5:I:186:DT:OP1	1:K:47:ALA:N	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:62:ILE:HG22	2:L:33:ALA:HB1	1.84	0.58
2:B:22:LEU:O	2:B:25:ASN:ND2	2.37	0.58
1:E:119:ILE:HG12	2:F:43:VAL:HG21	1.86	0.58
6:J:40:DG:OP1	4:N:86:ARG:HG3	2.03	0.58
1:A:63:ARG:NH1	6:J:194:DA:O3'	2.37	0.58
5:I:48:DC:H2''	5:I:49:DC:C5	2.39	0.58
3:G:20:ARG:NH2	6:J:135:DT:OP1	2.37	0.58
3:G:84:GLN:NE2	3:G:106:GLY:O	2.37	0.57
5:I:196:DC:H2''	5:I:197:DG:C8	2.40	0.57
1:A:126:LEU:HD22	1:E:113:HIS:CG	2.39	0.57
6:J:51:DC:H5''	1:K:83:ARG:HD2	1.86	0.57
5:I:117:DA:H4'	5:I:118:DT:OP1	2.05	0.56
2:F:84:MET:HE2	2:F:84:MET:HA	1.87	0.56
5:I:212:DT:H2''	5:I:213:DA:N7	2.21	0.56
1:K:97:GLU:O	1:K:101:VAL:HG13	2.06	0.56
5:I:222:DT:H2''	5:I:223:DG:N7	2.20	0.56
1:E:63:ARG:HH21	2:F:33:ALA:HB2	1.70	0.56
1:K:50:GLU:OE1	2:L:39:ARG:NE	2.32	0.56
3:M:29:ARG:NH1	4:N:35:GLU:OE1	2.39	0.56
1:E:103:LEU:O	1:E:107:THR:HG23	2.05	0.55
6:J:99:DG:H2''	6:J:100:DG:N7	2.21	0.55
5:I:185:DG:P	2:L:35:ARG:HH12	2.28	0.55
4:H:105:GLU:N	4:H:105:GLU:OE1	2.34	0.55
5:I:143:DA:H2''	5:I:144:DG:H5''	1.86	0.55
6:J:165:DC:H2''	6:J:166:DG:C8	2.41	0.55
1:K:57:SER:C	2:L:40:ARG:HH12	2.08	0.55
1:E:120:MET:HG3	1:E:122:LYS:HG2	1.87	0.55
5:I:49:DC:H2''	5:I:50:DT:H71	1.88	0.55
4:D:90:THR:HG22	4:D:93:GLU:OE1	2.06	0.55
5:I:248:DG:H2''	5:I:249:DA:C8	2.40	0.55
6:J:239:DT:H2''	6:J:240:DA:OP2	2.07	0.55
6:J:34:DG:H1'	6:J:35:DT:H5'	1.87	0.55
5:I:91:DA:H2''	5:I:92:DG:C8	2.42	0.55
2:B:92:ARG:HH21	4:D:101:LEU:HA	1.72	0.54
5:I:19:DA:H1'	5:I:20:DC:H5'	1.89	0.54
2:L:73:THR:HG21	2:L:81:VAL:HA	1.89	0.54
4:D:108:LYS:O	4:D:111:VAL:HG12	2.07	0.54
3:G:79:ILE:HG12	3:G:82:HIS:CE1	2.42	0.54
4:N:108:LYS:O	4:N:111:VAL:HG12	2.07	0.54
1:O:107:THR:HG23	1:O:123:ASP:HB2	1.90	0.54
5:I:180:DC:H2''	5:I:181:DG:C8	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:J:2:DT:H2'	6:J:3:DC:C6	2.42	0.54
6:J:3:DC:H2''	6:J:4:DG:C8	2.42	0.54
1:E:73:GLU:HG3	2:F:23:ARG:HB3	1.90	0.54
6:J:232:DT:H2''	6:J:233:DC:C5	2.43	0.54
1:K:83:ARG:HB2	2:L:80:THR:HG22	1.88	0.54
2:B:51:TYR:O	2:B:55:ARG:HG3	2.08	0.54
3:C:17:ARG:NH1	5:I:31:DA:OP2	2.41	0.54
5:I:184:DC:H5''	5:I:184:DC:H6	1.74	0.53
5:I:244:DT:H4'	5:I:245:DC:OP1	2.07	0.53
6:J:53:DC:H2''	6:J:54:DC:H5''	1.90	0.53
1:A:107:THR:HG23	1:A:123:ASP:HB2	1.91	0.53
3:C:62:ILE:HD11	4:D:62:MET:SD	2.48	0.53
6:J:113:DA:H1'	6:J:114:DG:H5'	1.91	0.53
5:I:193:DA:H1'	5:I:194:DA:C8	2.44	0.53
5:I:212:DT:H2''	5:I:213:DA:C8	2.44	0.53
6:J:152:DA:H1'	6:J:153:DG:C8	2.43	0.53
3:M:26:PRO:O	3:M:30:VAL:HG23	2.09	0.53
3:G:39:TYR:HB3	4:H:78:SER:HB2	1.90	0.53
5:I:193:DA:H1'	5:I:194:DA:N7	2.23	0.53
6:J:215:DT:H2''	6:J:216:DA:C8	2.44	0.53
1:K:121:PRO:HA	1:K:124:ILE:HG22	1.90	0.53
2:B:38:ALA:HB1	2:B:43:VAL:HG21	1.90	0.53
2:B:82:THR:HG23	2:B:84:MET:H	1.74	0.53
3:C:17:ARG:HG3	3:C:17:ARG:HH11	1.74	0.53
5:I:40:DA:H2''	5:I:41:DG:C8	2.44	0.53
6:J:216:DA:H2''	6:J:217:DG:H8	1.73	0.53
6:J:227:DA:H2''	6:J:228:DC:H5''	1.91	0.53
6:J:145:DC:H2''	6:J:146:DA:C8	2.44	0.53
1:A:61:LEU:O	2:B:36:ARG:NH1	2.42	0.52
2:F:78:ARG:NH1	2:F:80:THR:O	2.42	0.52
5:I:205:DA:H2''	5:I:206:DG:C8	2.44	0.52
2:P:91:LYS:HG3	2:P:96:THR:HG22	1.91	0.52
1:A:62:ILE:HD12	1:A:97:GLU:HG2	1.92	0.52
5:I:95:DG:H2''	5:I:96:DT:H6	1.74	0.52
6:J:165:DC:H2''	6:J:166:DG:H8	1.74	0.52
1:A:70:LEU:HD22	2:B:29:ILE:HD11	1.92	0.52
6:J:157:DC:H2''	6:J:158:DG:C8	2.44	0.52
1:E:107:THR:HG22	1:E:124:ILE:HA	1.91	0.52
5:I:186:DT:H2''	5:I:187:DG:C8	2.44	0.52
1:K:57:SER:O	2:L:40:ARG:NH1	2.37	0.52
5:I:197:DG:H2''	5:I:198:DG:H5''	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:61:LEU:HD21	2:L:40:ARG:HH21	1.75	0.52
3:G:79:ILE:H	3:G:82:HIS:HD1	1.58	0.52
5:I:184:DC:H2"	5:I:185:DG:C8	2.44	0.52
6:J:72:DC:H2"	6:J:73:DG:C8	2.44	0.52
1:K:65:LEU:HG	1:K:66:PRO:HD3	1.91	0.52
5:I:31:DA:H2"	5:I:32:DG:H5"	1.92	0.52
1:E:72:ARG:HH22	6:J:154:DC:P	2.32	0.52
1:O:54:TYR:CZ	2:P:36:ARG:HD3	2.45	0.52
2:B:35:ARG:NH1	6:J:185:DC:OP2	2.36	0.51
5:I:151:DC:H2"	5:I:152:DC:C6	2.45	0.51
5:I:247:DC:H2"	5:I:248:DG:H8	1.75	0.51
6:J:196:DC:H2"	6:J:197:DG:C8	2.46	0.51
6:J:226:DC:H2"	6:J:227:DA:H8	1.75	0.51
4:D:68:ASP:OD2	2:F:98:TYR:OH	2.26	0.51
6:J:123:DC:H2"	6:J:124:DT:H72	1.93	0.51
1:K:97:GLU:OE2	2:L:37:LEU:HD21	2.11	0.51
3:C:58:LEU:HD11	4:D:102:LEU:HD21	1.91	0.51
3:G:24:GLN:HB3	4:H:43:LYS:HE2	1.91	0.51
4:H:90:THR:HG22	4:H:93:GLU:OE1	2.09	0.51
5:I:178:DC:H2"	5:I:179:DG:C8	2.46	0.51
6:J:70:DC:H2"	6:J:71:DG:C8	2.46	0.51
1:O:57:SER:OG	2:P:40:ARG:NH2	2.42	0.51
2:F:71:THR:HG22	4:H:99:ARG:HH22	1.76	0.51
6:J:216:DA:H2"	6:J:217:DG:C8	2.46	0.51
2:F:92:ARG:HH21	4:H:101:LEU:HD23	1.76	0.51
5:I:52:DG:H2"	5:I:53:DG:H8	1.76	0.51
5:I:52:DG:H2"	5:I:53:DG:C8	2.46	0.50
5:I:3:DC:H5"	5:I:3:DC:H6	1.77	0.50
2:F:36:ARG:NH2	6:J:164:DA:OP1	2.26	0.50
6:J:192:DT:H5'	6:J:192:DT:C6	2.47	0.50
3:G:77:ARG:NH1	6:J:123:DC:OP1	2.44	0.50
6:J:192:DT:H6	6:J:192:DT:H5'	1.75	0.50
6:J:4:DG:H1'	6:J:5:DA:H5'	1.94	0.50
1:K:61:LEU:HD21	2:L:40:ARG:HE	1.75	0.50
5:I:215:DG:H5"	3:M:44:GLY:HA2	1.94	0.50
5:I:183:DA:H1'	5:I:184:DC:C5	2.47	0.50
5:I:226:DC:H2"	5:I:227:DG:C8	2.47	0.50
6:J:17:DC:H2"	6:J:18:DC:C5	2.46	0.50
5:I:234:DG:H2"	5:I:235:DC:OP2	2.12	0.50
1:K:89:VAL:HA	1:K:92:LEU:HD23	1.93	0.50
1:A:48:LEU:O	1:A:52:ARG:HG3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:29:DG:H2''	5:I:30:DG:C8	2.47	0.49
5:I:41:DG:H2'	5:I:42:DT:H71	1.93	0.49
6:J:33:DG:H5'	6:J:33:DG:C8	2.46	0.49
3:M:39:TYR:HB3	4:N:78:SER:HB2	1.93	0.49
1:O:119:ILE:HG22	2:P:50:ILE:CD1	2.42	0.49
5:I:164:DA:H2''	5:I:165:DA:H5'	1.93	0.49
5:I:172:DG:H2''	5:I:173:DG:C8	2.47	0.49
1:A:63:ARG:HB2	1:A:66:PRO:HG2	1.93	0.49
2:F:80:THR:HG23	5:I:102:DA:OP1	2.12	0.49
1:O:97:GLU:O	1:O:101:VAL:HG13	2.12	0.49
3:G:58:LEU:O	3:G:62:ILE:HG13	2.12	0.49
5:I:171:DG:C8	5:I:171:DG:H5'	2.47	0.49
5:I:3:DC:H2''	5:I:4:DG:C8	2.47	0.49
5:I:50:DT:H1'	5:I:51:DT:H5'	1.93	0.49
6:J:226:DC:H2''	6:J:227:DA:C8	2.47	0.49
4:H:108:LYS:HG2	4:H:109:HIS:HD1	1.78	0.49
5:I:207:DC:H2'	5:I:208:DT:H72	1.94	0.49
1:E:63:ARG:HH21	2:F:33:ALA:CB	2.25	0.49
5:I:139:DA:H2''	5:I:140:DG:C8	2.47	0.49
1:O:62:ILE:HD11	1:O:67:PHE:CD1	2.48	0.49
3:C:32:ARG:NH2	4:D:35:GLU:OE1	2.46	0.49
2:F:92:ARG:NH2	4:H:101:LEU:HD23	2.28	0.49
4:H:36:SER:HG	4:H:38:SER:HG	1.61	0.49
3:M:17:ARG:HH21	3:M:28:GLY:HA2	1.78	0.49
6:J:150:DC:H2''	6:J:151:DT:H71	1.94	0.49
1:K:119:ILE:O	2:L:47:SER:HB3	2.13	0.49
6:J:12:DC:H2''	6:J:13:DG:C8	2.47	0.49
6:J:213:DC:H1'	6:J:214:DC:H5'	1.93	0.48
2:L:84:MET:HB3	4:N:83:TYR:HE1	1.78	0.48
1:E:70:LEU:O	1:E:74:ILE:HG22	2.12	0.48
3:G:62:ILE:HD12	3:G:63:LEU:N	2.29	0.48
1:K:128:ARG:HH22	1:K:134:ARG:HE	1.61	0.48
5:I:241:DG:H2''	5:I:242:DA:C8	2.49	0.48
1:O:116:ARG:NH2	1:O:120:MET:HG2	2.28	0.48
2:P:49:LEU:HD23	2:P:49:LEU:H	1.79	0.48
5:I:62:DA:H1'	5:I:63:DC:H5'	1.96	0.48
5:I:67:DG:H2''	5:I:68:DG:C8	2.48	0.48
6:J:9:DT:H2''	6:J:10:DC:C6	2.47	0.48
6:J:196:DC:H2''	6:J:197:DG:H8	1.79	0.48
2:L:73:THR:HG21	2:L:81:VAL:HG22	1.96	0.48
5:I:208:DT:H2''	5:I:209:DG:H8	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:ASN:HB2	2:B:43:VAL:HG12	1.96	0.48
2:P:22:LEU:HB3	2:P:25:ASN:HD21	1.78	0.48
2:B:93:GLN:NE2	2:B:95:ARG:HD3	2.29	0.48
3:G:18:SER:OG	3:G:25:PHE:O	2.28	0.48
2:L:83:ALA:O	2:L:86:VAL:HG22	2.14	0.48
2:F:83:ALA:O	2:F:87:VAL:HG13	2.13	0.48
4:H:95:GLN:HG3	4:H:111:VAL:HG23	1.96	0.48
6:J:54:DC:H2'	6:J:55:DG:C8	2.48	0.48
1:K:62:ILE:HG12	1:K:97:GLU:OE1	2.14	0.48
2:L:35:ARG:HD2	2:L:46:ILE:HD12	1.96	0.48
2:P:43:VAL:CG1	2:P:46:ILE:HG22	2.44	0.48
1:E:48:LEU:O	1:E:51:ILE:HG12	2.13	0.47
4:H:62:MET:O	4:H:66:VAL:HG13	2.14	0.47
5:I:102:DA:H1'	5:I:103:DG:H5'	1.96	0.47
2:P:78:ARG:NH2	2:P:82:THR:HG22	2.29	0.47
1:A:97:GLU:O	1:A:101:VAL:HG13	2.14	0.47
3:C:63:LEU:HD22	4:D:45:LEU:HD13	1.96	0.47
2:F:31:LYS:HG3	2:F:51:TYR:CE1	2.49	0.47
3:G:27:VAL:HG11	3:G:49:VAL:HG12	1.95	0.47
6:J:142:DA:H2''	6:J:143:DG:C8	2.50	0.47
3:M:58:LEU:HD21	4:N:102:LEU:HD21	1.96	0.47
5:I:198:DG:H2''	5:I:199:DT:H5'	1.96	0.47
6:J:228:DC:H2''	6:J:229:DG:C8	2.50	0.47
6:J:53:DC:C2	6:J:54:DC:C5	3.02	0.47
3:C:79:ILE:HG12	3:C:82:HIS:CE1	2.49	0.47
5:I:248:DG:H2''	5:I:249:DA:H8	1.79	0.47
6:J:191:DT:H1'	6:J:192:DT:H5''	1.97	0.47
1:K:88:ALA:O	1:K:92:LEU:HD23	2.13	0.47
5:I:205:DA:OP1	2:L:80:THR:HG23	2.15	0.47
3:G:29:ARG:NH1	4:H:35:GLU:OE1	2.48	0.47
5:I:159:DG:H3'	5:I:160:DT:H71	1.96	0.47
1:E:63:ARG:HH11	1:E:66:PRO:HG2	1.80	0.47
1:E:79:LYS:HD3	1:E:82:LEU:HD11	1.95	0.47
5:I:109:DT:H2''	5:I:110:DA:C8	2.49	0.47
5:I:165:DA:H1'	5:I:166:DC:H5'	1.97	0.47
6:J:173:DC:H2''	6:J:174:DG:C8	2.49	0.47
6:J:147:DG:H2''	6:J:148:DC:C5	2.50	0.47
1:K:64:LYS:HG2	1:K:90:MET:HE1	1.96	0.47
6:J:245:DT:H2'	6:J:246:DC:C5	2.50	0.46
2:B:54:THR:O	2:B:57:VAL:HG22	2.15	0.46
4:D:95:GLN:O	4:D:98:VAL:HG12	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:243:DT:H2''	5:I:244:DT:OP2	2.16	0.46
1:A:122:LYS:HB3	1:E:113:HIS:HE1	1.79	0.46
2:F:30:THR:HB	2:F:32:PRO:HD2	1.97	0.46
5:I:166:DC:H2''	5:I:167:DG:C8	2.49	0.46
6:J:122:DG:H2''	6:J:123:DC:C5	2.50	0.46
4:H:114:GLY:O	4:H:118:VAL:HG23	2.16	0.46
6:J:21:DG:OP1	4:N:53:GLY:HA3	2.16	0.46
1:K:78:PHE:CD1	2:L:67:ARG:HG3	2.51	0.46
1:A:109:LEU:HD21	1:E:129:ARG:CZ	2.45	0.46
5:I:240:DG:H2''	5:I:241:DG:C8	2.51	0.46
5:I:124:DT:H2''	5:I:125:DC:C6	2.51	0.46
2:L:75:HIS:HB2	4:N:96:THR:HG21	1.97	0.46
4:H:54:ILE:HG22	4:H:55:SER:O	2.15	0.46
5:I:160:DT:H2''	5:I:161:DT:C6	2.51	0.46
6:J:192:DT:H2''	6:J:193:DA:C8	2.51	0.46
5:I:23:DG:C8	5:I:24:DT:H72	2.50	0.46
3:G:80:PRO:HD3	4:H:58:ALA:HB2	1.97	0.46
5:I:147:DA:H5'	5:I:147:DA:C8	2.51	0.46
3:G:17:ARG:HD3	6:J:134:DT:OP2	2.16	0.46
1:K:68:GLN:HG2	1:K:89:VAL:HG11	1.98	0.46
2:P:52:GLU:OE1	2:P:55:ARG:NH1	2.49	0.46
6:J:130:DT:OP2	6:J:130:DT:H2'	2.15	0.45
6:J:132:DA:H2''	6:J:133:DA:C8	2.51	0.45
6:J:19:DG:H2''	6:J:20:DA:N7	2.31	0.45
1:O:47:ALA:HB1	2:P:39:ARG:NH1	2.30	0.45
2:B:82:THR:HG23	2:B:84:MET:N	2.31	0.45
4:H:36:SER:HB2	4:H:63:ASN:HD22	1.81	0.45
5:I:27:5CM:H2''	5:I:28:DT:C7	2.46	0.45
6:J:241:DT:H2''	6:J:242:DA:OP2	2.16	0.45
2:L:62:LEU:HA	2:L:62:LEU:HD23	1.79	0.45
1:A:100:LEU:HD11	2:B:58:LEU:HD22	1.98	0.45
5:I:219:DA:H2''	5:I:220:DA:C8	2.51	0.45
6:J:217:DG:H2''	6:J:218:DT:H71	1.97	0.45
6:J:35:DT:H1'	6:J:36:DC:H5'	1.98	0.45
6:J:67:DG:H1'	6:J:68:DT:H5'	1.97	0.45
4:N:90:THR:HG22	4:N:91:SER:H	1.81	0.45
3:C:40:SER:OG	3:C:41:GLU:N	2.49	0.45
5:I:144:DG:H1'	5:I:145:DT:H5'	1.98	0.45
5:I:77:DC:H2''	5:I:78:DG:C8	2.51	0.45
2:L:72:TYR:HB3	2:L:85:ASP:OD1	2.16	0.45
2:B:38:ALA:O	2:B:43:VAL:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:244:DT:H2''	5:I:245:DC:C6	2.52	0.45
3:C:77:ARG:NH1	5:I:20:DC:H5''	2.31	0.45
1:E:60:LEU:HD22	1:E:93:GLN:HG2	1.98	0.45
3:G:37:GLY:HA3	3:G:39:TYR:CE2	2.52	0.45
5:I:12:DA:C8	5:I:13:DT:H72	2.52	0.45
6:J:155:DA:H2''	6:J:156:DC:H6	1.82	0.45
6:J:1:DA:N9	6:J:2:DT:H72	2.31	0.45
6:J:228:DC:H2''	6:J:229:DG:H8	1.81	0.45
5:I:11:DT:H2''	5:I:12:DA:OP2	2.16	0.45
5:I:161:DT:H1'	5:I:162:DA:C8	2.52	0.45
5:I:31:DA:H2''	5:I:32:DG:H8	1.81	0.45
6:J:1:DA:P	6:J:1:DA:H3'	2.57	0.45
2:B:38:ALA:HB1	2:B:43:VAL:CG2	2.46	0.45
1:A:74:ILE:HD12	2:B:62:LEU:HD22	1.98	0.45
5:I:54:DC:H2''	5:I:55:DG:C8	2.52	0.45
1:E:69:ARG:HH22	5:I:91:DA:P	2.40	0.45
6:J:133:DA:H2'	6:J:134:DT:H71	1.98	0.45
2:L:66:ILE:O	2:L:70:VAL:HG13	2.17	0.45
1:E:48:LEU:HA	1:E:51:ILE:HG23	1.98	0.45
6:J:160:DT:H4'	6:J:161:DT:OP1	2.17	0.45
6:J:212:DC:H2''	6:J:213:DC:C5	2.51	0.45
5:I:164:DA:H3'	2:P:30:THR:OG1	2.16	0.45
3:G:68:ASN:O	3:G:71:ARG:HG2	2.17	0.44
5:I:230:DC:H1'	5:I:231:DT:H5'	1.99	0.44
2:B:33:ALA:HA	2:B:36:ARG:NH1	2.32	0.44
3:C:17:ARG:HD2	4:D:121:TYR:CE2	2.52	0.44
1:E:131:ARG:NH1	1:E:133:GLU:OE1	2.49	0.44
6:J:123:DC:H2''	6:J:124:DT:C7	2.47	0.44
6:J:82:DC:OP2	2:P:35:ARG:NH1	2.37	0.44
3:G:102:ILE:HG23	4:H:61:ILE:HD12	1.99	0.44
5:I:116:DA:H2''	5:I:117:DA:OP2	2.16	0.44
5:I:194:DA:H2''	5:I:195:DG:C8	2.52	0.44
5:I:216:DA:H1'	5:I:217:DC:H5'	1.98	0.44
6:J:134:DT:C2'	6:J:135:DT:H71	2.47	0.44
6:J:193:DA:H1'	6:J:194:DA:C8	2.52	0.44
2:L:59:LYS:HG2	2:L:63:GLU:OE2	2.17	0.44
3:G:77:ARG:HB2	3:G:77:ARG:HE	1.55	0.44
1:K:101:VAL:HG11	2:L:40:ARG:HG2	1.98	0.44
1:E:48:LEU:O	1:E:52:ARG:HG3	2.17	0.44
5:I:123:DC:C6	5:I:124:DT:H72	2.52	0.44
5:I:181:DG:H2''	5:I:182:DT:C7	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:18:DG:H1'	5:I:19:DA:C5	2.53	0.44
5:I:22:DC:H2''	5:I:23:DG:C8	2.53	0.44
3:M:63:LEU:HA	3:M:63:LEU:HD23	1.84	0.44
4:N:95:GLN:O	4:N:99:ARG:HG3	2.18	0.44
1:A:63:ARG:NH2	5:I:60:DA:H4'	2.33	0.44
3:C:26:PRO:O	3:C:30:VAL:HG22	2.18	0.44
3:C:102:ILE:HG23	4:D:61:ILE:HD13	1.99	0.44
1:E:128:ARG:HA	1:E:131:ARG:HG2	2.00	0.44
1:E:97:GLU:O	1:E:101:VAL:HG13	2.17	0.44
6:J:176:DG:H2''	6:J:177:DC:C6	2.52	0.44
6:J:82:DC:P	2:P:35:ARG:HH22	2.40	0.44
2:L:57:VAL:O	2:L:60:VAL:HG22	2.17	0.44
4:N:68:ASP:OD2	2:P:98:TYR:OH	2.31	0.44
1:O:99:TYR:OH	1:O:133:GLU:OE1	2.22	0.44
2:B:73:THR:HB	2:B:85:ASP:OD2	2.18	0.44
2:B:92:ARG:NH2	4:D:101:LEU:HA	2.32	0.44
5:I:136:DA:H2'	5:I:137:DC:C5	2.52	0.44
1:O:121:PRO:HD3	2:P:50:ILE:HG13	1.99	0.44
1:O:116:ARG:NH2	1:O:123:ASP:OD2	2.50	0.44
1:O:78:PHE:CZ	2:P:67:ARG:HB3	2.53	0.44
1:A:119:ILE:HD11	2:B:46:ILE:HG12	1.99	0.44
1:E:79:LYS:CB	1:E:82:LEU:HD21	2.44	0.44
1:E:74:ILE:HD11	2:F:63:GLU:HA	2.00	0.44
5:I:12:DA:H2'	5:I:12:DA:OP2	2.17	0.44
6:J:9:DT:H2''	6:J:10:DC:H6	1.83	0.44
1:K:126:LEU:HD22	1:O:113:HIS:CG	2.52	0.44
1:K:89:VAL:HA	1:K:92:LEU:CD2	2.48	0.44
2:L:26:ILE:HD12	2:L:59:LYS:HD2	2.00	0.44
1:E:45:THR:O	1:E:48:LEU:HD23	2.18	0.43
3:G:88:ARG:NH2	3:G:100:VAL:O	2.51	0.43
6:J:156:DC:C2	6:J:157:DC:C5	3.06	0.43
6:J:94:DG:C4'	6:J:95:DC:H5'	2.48	0.43
1:K:51:ILE:HG23	2:L:42:GLY:HA2	1.99	0.43
1:E:102:GLY:O	1:E:105:GLU:HG2	2.17	0.43
3:G:16:THR:O	3:G:19:SER:OG	2.29	0.43
5:I:149:DC:C6	5:I:149:DC:H5'	2.54	0.43
2:L:67:ARG:O	2:L:70:VAL:HG22	2.18	0.43
2:P:26:ILE:HD13	2:P:59:LYS:HG3	2.00	0.43
2:F:26:ILE:HD12	2:F:26:ILE:H	1.82	0.43
2:B:30:THR:HG21	5:I:61:DA:H5''	2.00	0.43
1:K:75:ALA:HB1	1:K:82:LEU:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:63:LEU:HA	3:G:63:LEU:HD12	1.83	0.43
5:I:95:DG:H2''	5:I:96:DT:C6	2.52	0.43
6:J:70:DC:H5''	2:L:45:ARG:CD	2.48	0.43
2:L:92:ARG:NH2	4:N:101:LEU:HD23	2.33	0.43
2:B:73:THR:HG21	2:B:81:VAL:HG22	2.00	0.43
1:E:65:LEU:HG	5:I:92:DG:OP2	2.19	0.43
1:E:73:GLU:HB2	2:F:23:ARG:HD3	2.00	0.43
2:F:50:ILE:HG23	2:F:54:THR:HG23	2.00	0.43
3:G:117:PRO:HG3	5:I:145:DT:OP2	2.18	0.43
3:G:31:HIS:ND1	3:G:48:PRO:HG3	2.33	0.43
5:I:188:DC:H2''	5:I:189:DG:C8	2.53	0.43
5:I:44:DA:H2''	5:I:45:DT:OP2	2.16	0.43
3:M:51:LEU:O	3:M:54:VAL:HG22	2.18	0.43
2:P:90:LEU:HB3	2:P:95:ARG:O	2.18	0.43
2:B:31:LYS:HG3	2:B:51:TYR:CE1	2.54	0.43
5:I:140:DG:H1'	5:I:141:DG:C8	2.53	0.43
5:I:91:DA:H2''	5:I:92:DG:H8	1.83	0.43
1:E:109:LEU:HD23	1:E:109:LEU:HA	1.70	0.43
5:I:172:DG:H2''	5:I:173:DG:N7	2.33	0.43
6:J:95:DC:H2''	6:J:96:DC:OP2	2.18	0.43
1:K:63:ARG:HD3	1:K:63:ARG:HA	1.49	0.43
2:F:51:TYR:O	2:F:55:ARG:HG3	2.19	0.43
6:J:113:DA:H2''	6:J:114:DG:C8	2.54	0.43
1:E:108:ASN:O	1:E:112:ILE:HG23	2.19	0.43
1:E:74:ILE:HG21	1:E:74:ILE:HD13	1.74	0.43
5:I:55:DG:H2''	5:I:56:DG:C8	2.54	0.43
6:J:75:DT:H2''	6:J:76:DG:OP2	2.19	0.43
3:M:66:ALA:HB2	3:M:83:LEU:HD23	2.00	0.43
1:O:108:ASN:HA	1:O:119:ILE:HD11	2.01	0.43
1:E:46:VAL:HG23	1:E:49:ARG:CZ	2.49	0.42
5:I:138:DT:C2	5:I:139:DA:N7	2.87	0.42
5:I:69:DG:H4'	5:I:70:DG:OP1	2.18	0.42
6:J:143:DG:H1'	6:J:144:DA:C8	2.53	0.42
6:J:30:DA:C2'	6:J:31:DT:H71	2.48	0.42
2:P:43:VAL:HG11	2:P:46:ILE:HG22	2.00	0.42
6:J:1:DA:C4	6:J:2:DT:H72	2.53	0.42
1:K:74:ILE:HG21	2:L:66:ILE:HD12	2.01	0.42
1:A:60:LEU:HD12	1:A:64:LYS:HE2	2.01	0.42
6:J:141:DT:H2''	6:J:142:DA:N7	2.34	0.42
6:J:193:DA:H1'	6:J:194:DA:N7	2.35	0.42
6:J:231:DG:H2''	6:J:232:DT:OP2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:N:44:VAL:O	4:N:48:VAL:HG12	2.19	0.42
1:E:39:HIS:CG	5:I:7:DT:H5''	2.54	0.42
5:I:31:DA:H2''	5:I:32:DG:C8	2.55	0.42
3:M:29:ARG:NH2	4:N:40:TYR:HE2	2.16	0.42
1:A:131:ARG:NH1	1:A:133:GLU:OE2	2.49	0.42
5:I:184:DC:H2''	5:I:185:DG:H8	1.82	0.42
6:J:167:DC:H2''	6:J:168:DA:C8	2.55	0.42
2:P:33:ALA:HA	2:P:36:ARG:NH1	2.33	0.42
3:G:24:GLN:HG2	3:G:56:GLU:OE1	2.18	0.42
6:J:56:DC:H2''	6:J:57:DT:OP2	2.19	0.42
2:P:35:ARG:O	2:P:39:ARG:HG2	2.18	0.42
5:I:179:DG:OP2	5:I:179:DG:H2'	2.20	0.42
6:J:240:DA:H2''	6:J:241:DT:OP2	2.20	0.42
1:K:103:LEU:O	1:K:107:THR:OG1	2.29	0.42
1:K:128:ARG:NH2	1:K:134:ARG:HG3	2.34	0.42
2:L:30:THR:OG1	2:L:32:PRO:HD2	2.20	0.42
5:I:90:DA:H1'	5:I:91:DA:C8	2.55	0.42
6:J:154:DC:C2	6:J:155:DA:C8	3.08	0.42
6:J:202:DG:H2''	6:J:203:DG:C8	2.55	0.42
5:I:159:DG:H4'	5:I:160:DT:OP1	2.20	0.42
6:J:110:DC:H2''	6:J:111:DC:C6	2.55	0.42
1:O:115:LYS:HD3	1:O:115:LYS:HA	1.67	0.42
1:A:113:HIS:CG	1:E:126:LEU:HD22	2.55	0.41
2:B:34:ILE:CG2	2:B:54:THR:HG21	2.48	0.41
3:M:58:LEU:HB3	4:N:65:PHE:HZ	1.85	0.41
1:O:48:LEU:O	1:O:52:ARG:HG3	2.19	0.41
1:A:54:TYR:HA	1:A:57:SER:HB3	2.02	0.41
2:B:90:LEU:HB3	2:B:95:ARG:O	2.19	0.41
6:J:248:DG:H2''	6:J:249:DA:C8	2.55	0.41
6:J:32:DT:H2''	6:J:33:DG:C8	2.54	0.41
2:L:50:ILE:O	2:L:53:GLU:HB3	2.19	0.41
3:M:64:GLU:O	3:M:68:ASN:ND2	2.41	0.41
2:B:35:ARG:O	2:B:39:ARG:HG2	2.19	0.41
5:I:153:DT:H1'	5:I:154:DT:H5'	2.02	0.41
4:D:41:VAL:HA	4:D:44:VAL:HG12	2.02	0.41
5:I:6:DA:C2'	5:I:7:DT:H5'	2.49	0.41
1:A:129:ARG:NH2	1:E:106:ASP:OD1	2.54	0.41
5:I:23:DG:H2''	5:I:24:DT:O5'	2.19	0.41
6:J:154:DC:C2	6:J:155:DA:N7	2.89	0.41
2:L:52:GLU:O	2:L:55:ARG:HB2	2.20	0.41
2:B:57:VAL:O	2:B:60:VAL:HG12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:96:CYS:SG	2:F:62:LEU:HD13	2.60	0.41
2:F:35:ARG:O	2:F:39:ARG:HG2	2.21	0.41
2:F:45:ARG:NE	6:J:173:DC:H4'	2.35	0.41
5:I:75:DC:H2''	5:I:76:DG:OP2	2.21	0.41
6:J:136:DG:H1'	6:J:137:DG:C8	2.56	0.41
2:L:77:LYS:HE3	2:L:77:LYS:HB3	1.84	0.41
1:E:61:LEU:H	1:E:97:GLU:CD	2.24	0.41
5:I:22:DC:H2''	5:I:23:DG:H8	1.86	0.41
2:B:31:LYS:HG3	2:B:51:TYR:CD1	2.55	0.41
2:B:77:LYS:HE3	2:B:77:LYS:HB2	1.85	0.41
3:G:54:VAL:HG13	4:H:110:ALA:HB1	2.03	0.41
5:I:143:DA:C2'	5:I:144:DG:H5''	2.51	0.41
1:K:65:LEU:HA	1:K:68:GLN:HB2	2.03	0.41
1:K:66:PRO:O	1:K:69:ARG:HG3	2.20	0.41
6:J:31:DT:H5''	3:M:16:THR:HA	2.02	0.41
3:C:63:LEU:HD23	3:C:63:LEU:HA	1.95	0.41
3:G:29:ARG:HH21	4:H:40:TYR:HE1	1.67	0.41
5:I:205:DA:P	2:L:78:ARG:HG2	2.61	0.41
5:I:97:DG:O4'	5:I:97:DG:OP1	2.38	0.41
3:M:63:LEU:HD13	4:N:45:LEU:HB2	2.03	0.41
2:P:26:ILE:O	2:P:55:ARG:HD3	2.21	0.41
2:F:90:LEU:HB3	2:F:95:ARG:O	2.20	0.41
5:I:219:DA:H2''	5:I:220:DA:H8	1.86	0.41
6:J:13:DG:H8	6:J:13:DG:OP2	2.04	0.41
6:J:96:DC:H2''	6:J:97:DA:H5'	2.03	0.41
4:D:97:ALA:O	4:D:101:LEU:HD12	2.21	0.40
5:I:177:DG:H2''	5:I:178:DC:C6	2.56	0.40
5:I:68:DG:H1'	5:I:69:DG:C8	2.56	0.40
6:J:113:DA:H2''	6:J:114:DG:H8	1.86	0.40
1:K:71:VAL:HG23	2:L:66:ILE:HD13	2.02	0.40
1:A:119:ILE:HD13	2:B:43:VAL:HG11	2.02	0.40
1:E:60:LEU:HD13	1:E:93:GLN:CD	2.42	0.40
4:H:102:LEU:HA	4:H:102:LEU:HD23	1.89	0.40
5:I:17:DT:H4'	5:I:18:DG:OP1	2.22	0.40
6:J:31:DT:C2'	6:J:32:DT:H71	2.49	0.40
6:J:104:DT:OP1	4:N:32:SER:HA	2.21	0.40
5:I:227:DG:H5'	4:N:33:ARG:HG2	2.02	0.40
4:D:72:ARG:HB3	4:D:101:LEU:HD22	2.03	0.40
1:E:107:THR:CG2	1:E:124:ILE:HA	2.51	0.40
5:I:10:DA:H2''	5:I:11:DT:OP2	2.20	0.40
5:I:153:DT:C2'	5:I:154:DT:H71	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:VAL:HG13	2:B:66:ILE:HD11	2.03	0.40
2:B:96:THR:HB	3:G:100:VAL:HG12	2.04	0.40
3:C:116:LEU:HA	3:C:116:LEU:HD23	1.93	0.40
5:I:198:DG:H2'	5:I:199:DT:C6	2.57	0.40
5:I:240:DG:H2''	5:I:241:DG:H8	1.85	0.40
1:K:119:ILE:HG12	2:L:43:VAL:HG13	2.03	0.40
2:L:84:MET:O	2:L:88:TYR:HD1	2.04	0.40
3:M:40:SER:OG	3:M:41:GLU:N	2.54	0.40
3:C:32:ARG:HH22	4:D:35:GLU:CD	2.25	0.40
2:F:67:ARG:O	2:F:71:THR:HG23	2.21	0.40
4:H:66:VAL:HA	4:H:69:ILE:HG22	2.03	0.40
5:I:191:DT:H2''	5:I:192:DT:H72	2.04	0.40
5:I:78:DG:H2''	5:I:79:DT:C6	2.57	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	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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
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%)	83 (99%)	1 (1%)	71	87
1	E	85/113 (75%)	83 (98%)	2 (2%)	49	75
1	K	77/113 (68%)	75 (97%)	2 (3%)	46	73
1	O	84/113 (74%)	82 (98%)	2 (2%)	49	75
2	B	67/81 (83%)	65 (97%)	2 (3%)	41	70
2	F	65/81 (80%)	63 (97%)	2 (3%)	40	69
2	L	58/81 (72%)	55 (95%)	3 (5%)	23	53
2	P	67/81 (83%)	67 (100%)	0	100	100
3	C	81/102 (79%)	81 (100%)	0	100	100
3	G	85/102 (83%)	79 (93%)	6 (7%)	14	42
3	M	83/102 (81%)	81 (98%)	2 (2%)	49	75
4	D	79/107 (74%)	77 (98%)	2 (2%)	47	74
4	H	77/107 (72%)	73 (95%)	4 (5%)	23	53
4	N	79/107 (74%)	77 (98%)	2 (2%)	47	74
All	All	1071/1403 (76%)	1041 (97%)	30 (3%)	43	71

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	ARG
2	B	22	LEU
2	B	93	GLN
4	D	51	ASP
4	D	55	SER
1	E	63	ARG
1	E	131	ARG
2	F	59	LYS
2	F	92	ARG
3	G	9	LYS
3	G	15	LYS
3	G	17	ARG
3	G	71	ARG
3	G	83	LEU
3	G	99	ARG
4	H	51	ASP
4	H	67	ASN
4	H	80	LEU
4	H	108	LYS
1	K	49	ARG
1	K	134	ARG
2	L	25	ASN
2	L	52	GLU
2	L	92	ARG
3	M	74	LYS
3	M	81	ARG
4	N	34	LYS
4	N	51	ASP
1	O	48	LEU
1	O	72	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	73	ASN
3	G	73	ASN
1	K	85	GLN

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$
6	5CM	J	222	5,6	15,21,22	1.35	1 (6%)	19,30,33	1.44	4 (21%)
5	5CM	I	27	5,6	15,21,22	1.37	1 (6%)	19,30,33	1.46	4 (21%)
5	5CM	I	130	5,6	15,21,22	1.31	1 (6%)	19,30,33	1.46	4 (21%)

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
6	5CM	J	222	5,6	-	1/4/21/22	0/2/2/2
5	5CM	I	27	5,6	-	1/4/21/22	0/2/2/2
5	5CM	I	130	5,6	-	1/4/21/22	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	27	5CM	C5-C4	4.89	1.48	1.41
6	J	222	5CM	C5-C4	4.77	1.48	1.41
5	I	130	5CM	C5-C4	4.58	1.48	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	27	5CM	C2-N3-C4	3.58	120.33	116.02
5	I	130	5CM	C2-N3-C4	3.24	119.93	116.02
6	J	222	5CM	C2-N3-C4	3.24	119.93	116.02
6	J	222	5CM	C5-C6-N1	-3.05	118.91	122.19
5	I	130	5CM	C5-C6-N1	-2.91	119.05	122.19
5	I	27	5CM	C5-C6-N1	-2.66	119.32	122.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	130	5CM	N4-C4-N3	2.43	120.47	117.03
5	I	130	5CM	C5A-C5-C4	-2.39	119.31	121.72
6	J	222	5CM	N4-C4-N3	2.36	120.37	117.03
5	I	27	5CM	O3'-C3'-C2'	-2.15	103.20	110.90
5	I	27	5CM	N4-C4-N3	2.13	120.04	117.03
6	J	222	5CM	C5A-C5-C4	-2.05	119.65	121.72

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	J	222	5CM	O4'-C1'-N1-C6
5	I	27	5CM	O4'-C1'-N1-C6
5	I	130	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 carbohydrates 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 ⓘ

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(Å ²)	Q<0.9
1	A	96/139 (69%)	-0.33	0 100 100	48, 64, 84, 103	0
1	E	97/139 (69%)	-0.27	0 100 100	59, 74, 100, 115	0
1	K	88/139 (63%)	-0.26	0 100 100	61, 75, 103, 129	0
1	O	96/139 (69%)	-0.19	1 (1%) 82 70	58, 75, 100, 121	0
2	B	82/106 (77%)	-0.31	0 100 100	51, 64, 84, 92	0
2	F	79/106 (74%)	-0.22	1 (1%) 77 61	53, 73, 90, 111	0
2	L	69/106 (65%)	-0.28	0 100 100	55, 76, 91, 95	0
2	P	81/106 (76%)	-0.26	0 100 100	53, 73, 101, 109	0
3	C	102/133 (76%)	-0.19	1 (0%) 82 70	61, 73, 94, 103	0
3	G	109/133 (81%)	-0.23	1 (0%) 84 72	57, 76, 97, 118	0
3	M	106/133 (79%)	-0.30	0 100 100	55, 71, 94, 115	0
4	D	92/129 (71%)	-0.15	0 100 100	57, 75, 90, 106	0
4	H	91/129 (70%)	-0.25	0 100 100	63, 79, 94, 109	0
4	N	93/129 (72%)	-0.39	0 100 100	51, 72, 91, 118	0
5	I	243/250 (97%)	-0.61	0 100 100	78, 115, 145, 162	0
6	J	244/250 (97%)	-0.65	1 (0%) 92 86	74, 117, 145, 159	0
All	All	1768/2266 (78%)	-0.36	5 (0%) 94 89	48, 79, 133, 162	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	77	ARG	3.2
3	C	117	PRO	2.9
6	J	1	DA	2.8
2	F	101	GLY	2.0
1	O	81	ASP	2.0

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(\AA^2)	Q<0.9
5	5CM	I	130	20/21	0.78	0.21	98,124,138,154	0
6	5CM	J	222	20/21	0.90	0.18	99,113,121,123	0
5	5CM	I	27	20/21	0.93	0.17	93,111,123,132	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.