Full wwPDB X-ray Structure Validation Report ⓘ

Nov 24, 2021 – 01:10 pm GMT

PDB ID	:	7AZ5
Title	:	DNA polymerase sliding clamp from Escherichia coli with peptide 47 bound
Authors	:	Monsarrat, C.; Compain, G.; Andre, C.; Martiel, I.; Engilberge, S.; Olieric, V.; Wolff, P.; Brillet, K.; Landolfo, M.; Silva da Veiga, C.; Wagner, J.; Guichard, G.; Burnouf, D.Y.
Deposited on	:	2020-11-16
Resolution	:	1.87 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:

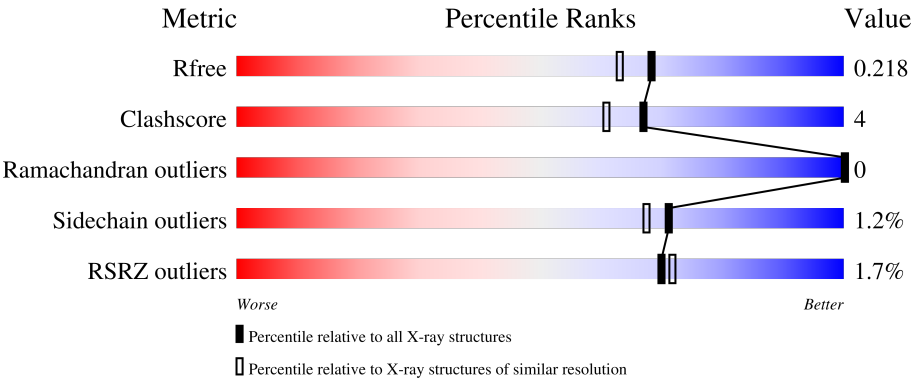
MolProbity	:	4.02b-467
Mogul	:	1.8.4 (270009), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 1.87 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	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	386	.% 87% 9% 5%
1	B	386	3% 87% 9% 5%
1	C	386	2% 84% 11% 5%
1	D	386	.% 89% 6% 5%
2	H	7	57% 43%

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Mol	Chain	Length	Quality of chain
2	I	7	57% 43%
2	J	7	86% 14%
2	K	7	86% 14%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12930 atoms, of which 30 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 Beta sliding clamp.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	2	0
			2878	1805	507	547	19			
1	B	368	Total	C	N	O	S	0	3	0
			2877	1805	505	548	19			
1	C	368	Total	C	N	O	S	0	3	0
			2883	1809	510	545	19			
1	D	368	Total	C	N	O	S	0	4	0
			2881	1809	508	545	19			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A073FMV0
A	-18	GLY	-	expression tag	UNP A0A073FMV0
A	-17	SER	-	expression tag	UNP A0A073FMV0
A	-16	SER	-	expression tag	UNP A0A073FMV0
A	-15	HIS	-	expression tag	UNP A0A073FMV0
A	-14	HIS	-	expression tag	UNP A0A073FMV0
A	-13	HIS	-	expression tag	UNP A0A073FMV0
A	-12	HIS	-	expression tag	UNP A0A073FMV0
A	-11	HIS	-	expression tag	UNP A0A073FMV0
A	-10	HIS	-	expression tag	UNP A0A073FMV0
A	-9	SER	-	expression tag	UNP A0A073FMV0
A	-8	SER	-	expression tag	UNP A0A073FMV0
A	-7	GLY	-	expression tag	UNP A0A073FMV0
A	-6	LEU	-	expression tag	UNP A0A073FMV0
A	-5	VAL	-	expression tag	UNP A0A073FMV0
A	-4	PRO	-	expression tag	UNP A0A073FMV0
A	-3	ARG	-	expression tag	UNP A0A073FMV0
A	-2	GLY	-	expression tag	UNP A0A073FMV0
A	-1	SER	-	expression tag	UNP A0A073FMV0
A	0	HIS	-	expression tag	UNP A0A073FMV0
B	-19	MET	-	initiating methionine	UNP A0A073FMV0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP A0A073FMV0
B	-17	SER	-	expression tag	UNP A0A073FMV0
B	-16	SER	-	expression tag	UNP A0A073FMV0
B	-15	HIS	-	expression tag	UNP A0A073FMV0
B	-14	HIS	-	expression tag	UNP A0A073FMV0
B	-13	HIS	-	expression tag	UNP A0A073FMV0
B	-12	HIS	-	expression tag	UNP A0A073FMV0
B	-11	HIS	-	expression tag	UNP A0A073FMV0
B	-10	HIS	-	expression tag	UNP A0A073FMV0
B	-9	SER	-	expression tag	UNP A0A073FMV0
B	-8	SER	-	expression tag	UNP A0A073FMV0
B	-7	GLY	-	expression tag	UNP A0A073FMV0
B	-6	LEU	-	expression tag	UNP A0A073FMV0
B	-5	VAL	-	expression tag	UNP A0A073FMV0
B	-4	PRO	-	expression tag	UNP A0A073FMV0
B	-3	ARG	-	expression tag	UNP A0A073FMV0
B	-2	GLY	-	expression tag	UNP A0A073FMV0
B	-1	SER	-	expression tag	UNP A0A073FMV0
B	0	HIS	-	expression tag	UNP A0A073FMV0
C	-19	MET	-	initiating methionine	UNP A0A073FMV0
C	-18	GLY	-	expression tag	UNP A0A073FMV0
C	-17	SER	-	expression tag	UNP A0A073FMV0
C	-16	SER	-	expression tag	UNP A0A073FMV0
C	-15	HIS	-	expression tag	UNP A0A073FMV0
C	-14	HIS	-	expression tag	UNP A0A073FMV0
C	-13	HIS	-	expression tag	UNP A0A073FMV0
C	-12	HIS	-	expression tag	UNP A0A073FMV0
C	-11	HIS	-	expression tag	UNP A0A073FMV0
C	-10	HIS	-	expression tag	UNP A0A073FMV0
C	-9	SER	-	expression tag	UNP A0A073FMV0
C	-8	SER	-	expression tag	UNP A0A073FMV0
C	-7	GLY	-	expression tag	UNP A0A073FMV0
C	-6	LEU	-	expression tag	UNP A0A073FMV0
C	-5	VAL	-	expression tag	UNP A0A073FMV0
C	-4	PRO	-	expression tag	UNP A0A073FMV0
C	-3	ARG	-	expression tag	UNP A0A073FMV0
C	-2	GLY	-	expression tag	UNP A0A073FMV0
C	-1	SER	-	expression tag	UNP A0A073FMV0
C	0	HIS	-	expression tag	UNP A0A073FMV0
D	-19	MET	-	initiating methionine	UNP A0A073FMV0
D	-18	GLY	-	expression tag	UNP A0A073FMV0
D	-17	SER	-	expression tag	UNP A0A073FMV0

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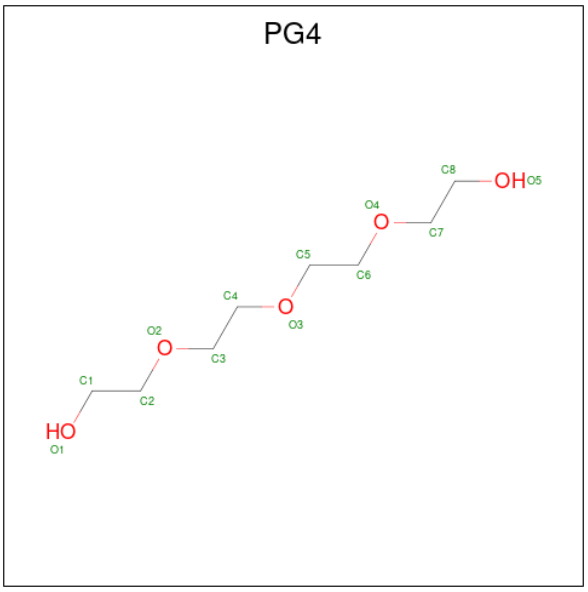
Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP A0A073FMV0
D	-15	HIS	-	expression tag	UNP A0A073FMV0
D	-14	HIS	-	expression tag	UNP A0A073FMV0
D	-13	HIS	-	expression tag	UNP A0A073FMV0
D	-12	HIS	-	expression tag	UNP A0A073FMV0
D	-11	HIS	-	expression tag	UNP A0A073FMV0
D	-10	HIS	-	expression tag	UNP A0A073FMV0
D	-9	SER	-	expression tag	UNP A0A073FMV0
D	-8	SER	-	expression tag	UNP A0A073FMV0
D	-7	GLY	-	expression tag	UNP A0A073FMV0
D	-6	LEU	-	expression tag	UNP A0A073FMV0
D	-5	VAL	-	expression tag	UNP A0A073FMV0
D	-4	PRO	-	expression tag	UNP A0A073FMV0
D	-3	ARG	-	expression tag	UNP A0A073FMV0
D	-2	GLY	-	expression tag	UNP A0A073FMV0
D	-1	SER	-	expression tag	UNP A0A073FMV0
D	0	HIS	-	expression tag	UNP A0A073FMV0

- Molecule 2 is a protein called Peptide 47.

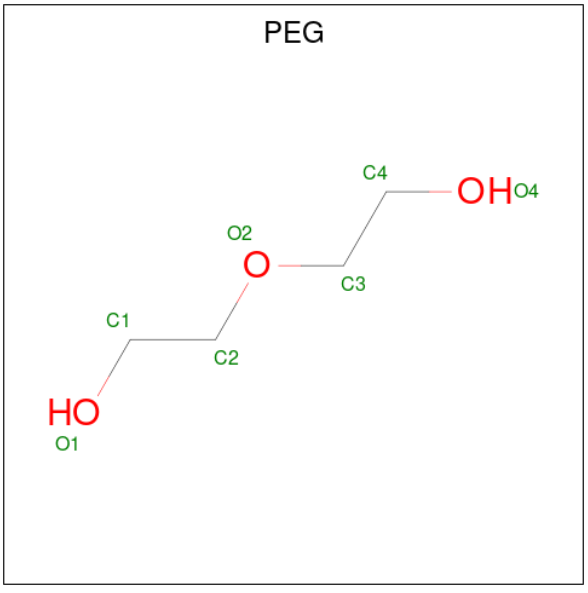
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	7	Total	C	N	O	0	0	0
			59	41	7	11			
2	I	7	Total	C	N	O	0	0	0
			59	41	7	11			
2	J	7	Total	C	N	O	0	0	0
			59	41	7	11			
2	K	7	Total	C	N	O	0	0	0
			59	41	7	11			

- Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	8	5		
3	B	1	Total	C	O	0	0
			13	8	5		
3	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			17	4	10	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C H O 17 4 10 3	0	0
4	C	1	Total C H O 17 4 10 3	0	0

- Molecule 5 is water.

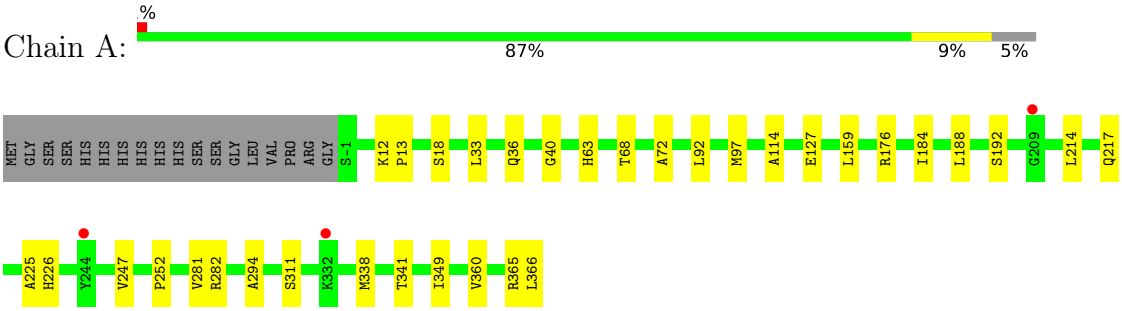
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	246	Total O 254 254	0	8
5	B	234	Total O 241 241	0	7
5	C	242	Total O 251 251	0	9
5	D	281	Total O 294 294	0	13
5	H	9	Total O 9 9	0	0
5	I	11	Total O 12 12	0	1
5	J	10	Total O 10 10	0	0
5	K	14	Total O 14 14	0	0

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.

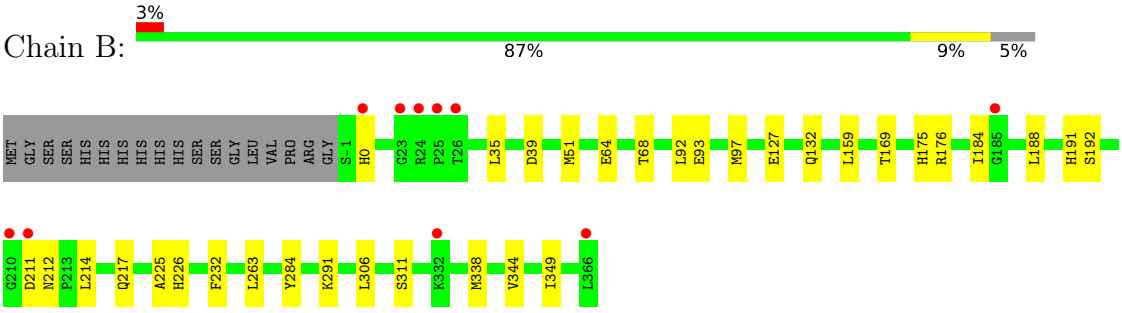
• Molecule 1: Beta sliding clamp

Chain A:



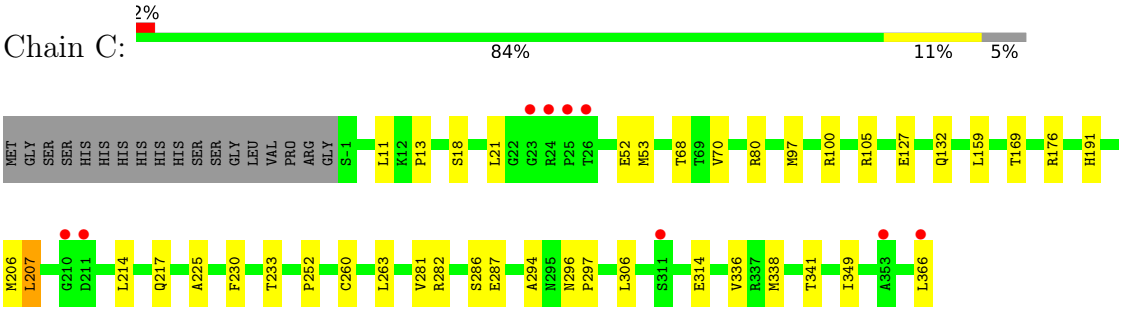
• Molecule 1: Beta sliding clamp

Chain B:



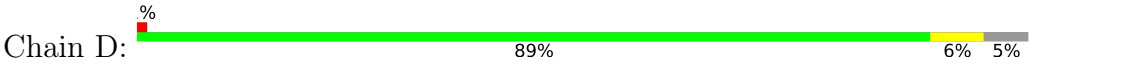
• Molecule 1: Beta sliding clamp

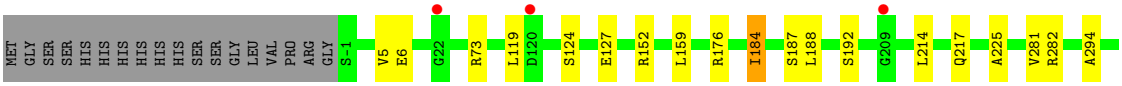
Chain C:

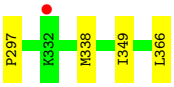


• Molecule 1: Beta sliding clamp

Chain D:





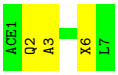


• Molecule 2: Peptide 47

Chain H:

• Molecule 2: Peptide 47

Chain I:



• Molecule 2: Peptide 47

Chain J:

• Molecule 2: Peptide 47

Chain K:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	68.89Å 80.17Å 83.37Å 62.79° 71.29° 81.66°	Depositor
Resolution (Å)	27.35 – 1.87 65.25 – 1.87	Depositor EDS
% Data completeness (in resolution range)	72.8 (27.35-1.87) 72.8 (65.25-1.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.68 (at 1.87Å)	Xtriage
Refinement program	BUSTER	Depositor
R, R_{free}	0.177 , 0.211 0.184 , 0.218	Depositor DCC
R_{free} test set	4338 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å^2)	25.8	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12930	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.83% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, PG4, ALC, OIC, ACE

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	0.46	0/2929	0.64	0/3964
1	B	0.47	0/2933	0.64	0/3970
1	C	0.47	0/2937	0.64	0/3974
1	D	0.47	0/2938	0.64	0/3976
2	H	0.55	0/32	0.72	0/38
2	I	0.49	0/32	0.76	0/38
2	J	0.56	0/32	0.76	0/38
2	K	0.52	0/32	0.69	0/38
All	All	0.47	0/11865	0.64	0/16036

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	1
2	I	0	1
2	J	0	1
2	K	0	1
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	6	OIC	Mainchain

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Mol	Chain	Res	Type	Group
2	I	6	OIC	Mainchain
2	J	6	OIC	Mainchain
2	K	6	OIC	Mainchain

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	2878	0	2884	26	0
1	B	2877	0	2891	23	0
1	C	2883	0	2895	29	0
1	D	2881	0	2893	16	0
2	H	59	0	64	2	0
2	I	59	0	64	2	0
2	J	59	0	64	0	0
2	K	59	0	64	0	0
3	A	13	0	18	3	0
3	B	13	0	18	0	0
3	D	13	0	18	2	0
4	A	7	10	10	3	0
4	B	7	10	10	0	0
4	C	7	10	10	1	0
5	A	254	0	0	0	0
5	B	241	0	0	2	0
5	C	251	0	0	2	0
5	D	294	0	0	1	0
5	H	9	0	0	0	0
5	I	12	0	0	0	0
5	J	10	0	0	0	0
5	K	14	0	0	0	0
All	All	12900	30	11903	92	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (92) 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:A:365:ARG:HH12	1:D:152:ARG:HG2	1.31	0.94
1:C:206:MET:HE2	1:C:230:PHE:HB3	1.49	0.92
1:A:184:ILE:HD11	1:A:188:LEU:HD11	1.57	0.87
1:B:68:THR:HG21	1:B:92:LEU:HD11	1.70	0.73
1:C:68:THR:CG2	1:C:97:MET:HE2	2.22	0.69
1:A:247:VAL:HG12	1:A:360:VAL:HG11	1.79	0.65
1:C:53:MET:SD	1:C:206:MET:HE3	2.38	0.64
1:C:206:MET:HE2	1:C:230:PHE:CB	2.27	0.63
1:A:365:ARG:HB2	2:H:3:ALC:HZ3	1.80	0.63
1:C:53:MET:SD	1:C:206:MET:CE	2.86	0.63
1:D:184:ILE:HD11	1:D:188:LEU:HD21	1.81	0.63
1:B:35:LEU:HD12	1:B:97:MET:CE	2.29	0.62
1:A:18:SER:HB3	3:A:401:PG4:C6	2.32	0.60
1:A:68:THR:HG21	1:A:92:LEU:HD21	1.81	0.60
1:A:18:SER:HB3	3:A:401:PG4:H61	1.83	0.60
1:A:247:VAL:HG13	2:H:5:LEU:HD22	1.83	0.60
1:A:114:ALA:HA	4:A:402:PEG:H11	1.85	0.59
1:B:0:HIS:CG	1:B:0:HIS:O	2.58	0.56
1:D:159:LEU:HD11	1:D:192:SER:HB2	1.86	0.56
1:A:159:LEU:HD11	1:A:192:SER:HB2	1.89	0.55
1:C:127:GLU:HG2	1:C:217:GLN:HG2	1.89	0.54
1:B:51:MET:HE3	1:B:232:PHE:HE1	1.74	0.53
1:B:127:GLU:HG2	1:B:217:GLN:HG2	1.90	0.53
1:A:214:LEU:HD11	1:A:225:ALA:HB1	1.89	0.53
1:D:127:GLU:HG2	1:D:217:GLN:HG2	1.91	0.53
1:C:53:MET:SD	1:C:206:MET:HE1	2.49	0.53
1:B:214:LEU:HD11	1:B:225:ALA:HB1	1.90	0.53
1:D:214:LEU:HD11	1:D:225:ALA:HB1	1.92	0.52
1:A:127:GLU:HG2	1:A:217:GLN:HG2	1.90	0.52
1:A:36:GLN:HE21	4:A:402:PEG:H41	1.74	0.52
1:C:191:HIS:HE1	5:C:694[B]:HOH:O	1.93	0.52
1:A:365:ARG:NH1	1:D:152:ARG:HG2	2.13	0.52
1:B:35:LEU:HD12	1:B:97:MET:HE3	1.91	0.51
1:B:344:VAL:HA	2:I:3:ALC:HE12	1.93	0.51
1:D:73:ARG:HB2	3:D:401:PG4:H32	1.93	0.51
1:A:68:THR:CG2	1:A:97:MET:HE2	2.41	0.50
1:B:35:LEU:HD12	1:B:97:MET:HE1	1.93	0.50
1:A:247:VAL:CG1	1:A:360:VAL:HG11	2.42	0.49
1:D:73:ARG:HH11	3:D:401:PG4:H12	1.78	0.48
1:B:284:TYR:HB2	1:B:291:LYS:HB3	1.95	0.48
1:B:184:ILE:HD11	1:B:188:LEU:HD11	1.94	0.48
1:D:338:MET:HG2	1:D:349:ILE:HG12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:HIS:CE1	5:C:694[B]:HOH:O	2.66	0.48
1:B:159:LEU:HD11	1:B:192:SER:HB2	1.95	0.48
1:D:124[A]:SER:OG	1:D:127:GLU:HG3	2.14	0.47
1:B:132:GLN:HE21	1:B:212:ASN:HB2	1.80	0.47
1:A:281:VAL:HG12	1:A:294:ALA:HB2	1.96	0.47
1:C:282:ARG:HG2	1:C:366:LEU:HD22	1.96	0.47
1:D:282:ARG:HG2	1:D:366:LEU:HD22	1.97	0.47
1:C:13:PRO:HA	1:C:230:PHE:HE2	1.81	0.46
1:C:132:GLN:HE21	1:C:207:LEU:HA	1.80	0.46
1:C:214:LEU:HD11	1:C:225:ALA:HB1	1.97	0.46
1:D:281:VAL:HG12	1:D:294:ALA:HB2	1.98	0.46
1:B:132:GLN:NE2	1:B:212:ASN:HB2	2.31	0.45
1:D:119:LEU:HD23	1:D:119:LEU:HA	1.85	0.45
1:A:252:PRO:HA	1:A:341:THR:HG22	1.99	0.45
1:A:338:MET:HG2	1:A:349:ILE:HG12	1.98	0.45
1:C:52:GLU:HG2	1:C:233:THR:HB	1.99	0.45
1:B:338:MET:HG2	1:B:349:ILE:HG12	1.99	0.44
1:C:68:THR:HG21	1:C:97:MET:HE2	1.98	0.44
1:B:0:HIS:NE2	1:B:64:GLU:HG3	2.33	0.44
1:B:191:HIS:CE1	5:B:541:HOH:O	2.71	0.44
1:A:18:SER:HB3	3:A:401:PG4:H62	2.00	0.44
1:C:70:VAL:HB	1:C:97:MET:HE1	2.00	0.43
1:A:33:LEU:HG	1:A:72:ALA:HB2	2.00	0.43
1:A:36:GLN:HE21	4:A:402:PEG:C4	2.31	0.43
1:C:80:ARG:HH12	4:C:401:PEG:H21	1.83	0.43
1:C:18:SER:HA	1:C:21:LEU:HD13	1.99	0.43
1:C:338:MET:HG2	1:C:349:ILE:HG12	2.01	0.43
1:B:159:LEU:O	1:B:169:THR:HA	2.19	0.43
1:C:263:LEU:HD12	1:C:306:LEU:HD22	2.00	0.43
1:C:260:CYS:HB2	1:C:336:VAL:HG23	2.01	0.43
1:B:68:THR:HB	1:B:92:LEU:HD21	2.01	0.43
1:B:175:HIS:CE1	2:I:2:GLN:HG2	2.53	0.43
1:C:286:SER:HA	1:C:314:GLU:HG2	2.00	0.43
1:C:68:THR:HG23	1:C:97:MET:HE2	2.00	0.42
1:C:100:ARG:HH21	1:C:105:ARG:HH12	1.67	0.42
1:B:191:HIS:HE1	5:B:541:HOH:O	2.01	0.42
1:A:40:GLY:HA2	1:A:63:HIS:NE2	2.34	0.42
1:A:68:THR:CG2	1:A:97:MET:CE	2.98	0.42
1:A:12:LYS:HB3	1:A:13:PRO:HD3	2.01	0.42
1:A:282:ARG:HE	1:A:366:LEU:HD22	1.84	0.41
1:C:252:PRO:HA	1:C:341:THR:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:ASN:HB2	1:C:297:PRO:CD	2.50	0.41
1:C:281:VAL:HG12	1:C:294:ALA:HB2	2.02	0.41
1:D:184:ILE:HG12	1:D:188:LEU:HD11	2.01	0.41
1:D:5:VAL:HG12	1:D:6:GLU:O	2.21	0.41
1:B:263:LEU:HD12	1:B:306:LEU:HD22	2.02	0.41
1:C:11:LEU:HD23	1:C:11:LEU:HA	1.89	0.41
1:C:159:LEU:O	1:C:169:THR:HA	2.21	0.40
1:D:297:PRO:HD2	5:D:527:HOH:O	2.21	0.40
1:B:51:MET:HE3	1:B:232:PHE:CE1	2.54	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	368/386 (95%)	361 (98%)	7 (2%)	0	100	100
1	B	369/386 (96%)	359 (97%)	10 (3%)	0	100	100
1	C	369/386 (96%)	359 (97%)	10 (3%)	0	100	100
1	D	370/386 (96%)	362 (98%)	8 (2%)	0	100	100
2	H	3/7 (43%)	3 (100%)	0	0	100	100
2	I	3/7 (43%)	3 (100%)	0	0	100	100
2	J	3/7 (43%)	3 (100%)	0	0	100	100
2	K	3/7 (43%)	3 (100%)	0	0	100	100
All	All	1488/1572 (95%)	1453 (98%)	35 (2%)	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	317/330 (96%)	314 (99%)	3 (1%)	78 76
1	B	318/330 (96%)	312 (98%)	6 (2%)	57 49
1	C	317/330 (96%)	314 (99%)	3 (1%)	78 76
1	D	317/330 (96%)	314 (99%)	3 (1%)	78 76
2	H	4/4 (100%)	4 (100%)	0	100 100
2	I	4/4 (100%)	4 (100%)	0	100 100
2	J	4/4 (100%)	4 (100%)	0	100 100
2	K	4/4 (100%)	4 (100%)	0	100 100
All	All	1285/1336 (96%)	1270 (99%)	15 (1%)	71 67

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	ARG
1	A	226	HIS
1	A	311	SER
1	B	39	ASP
1	B	93	GLU
1	B	176	ARG
1	B	211	ASP
1	B	226	HIS
1	B	311	SER
1	C	176	ARG
1	C	207	LEU
1	C	287	GLU
1	D	176	ARG
1	D	184	ILE
1	D	187	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	191	HIS
1	B	191	HIS
1	C	16	GLN
1	C	132	GLN
1	C	212	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

6 ligands are modelled in this entry.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	368/386 (95%)	-0.22	3 (0%) 86 87	19, 30, 56, 75	0
1	B	368/386 (95%)	-0.14	10 (2%) 54 56	17, 31, 59, 78	0
1	C	368/386 (95%)	-0.18	9 (2%) 59 60	16, 31, 59, 92	0
1	D	368/386 (95%)	-0.30	4 (1%) 80 82	16, 30, 52, 66	0
2	H	4/7 (57%)	-0.29	0 100 100	28, 31, 39, 41	0
2	I	4/7 (57%)	-0.74	0 100 100	24, 25, 26, 33	0
2	J	4/7 (57%)	-0.35	0 100 100	24, 27, 29, 38	0
2	K	4/7 (57%)	-0.36	0 100 100	25, 28, 33, 42	0
All	All	1488/1572 (94%)	-0.21	26 (1%) 70 72	16, 31, 58, 92	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	244	TYR	5.5
1	C	23	GLY	4.6
1	B	24	ARG	4.3
1	C	26	THR	4.2
1	D	209	GLY	3.9
1	C	24	ARG	3.8
1	B	0	HIS	3.5
1	B	366	LEU	3.4
1	C	311	SER	3.2
1	B	23	GLY	3.1
1	A	209	GLY	3.1
1	C	210	GLY	3.0
1	B	211	ASP	3.0
1	B	210	GLY	2.9
1	B	26	THR	2.9
1	D	120	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	332	LYS	2.7
1	D	332	LYS	2.6
1	B	332	LYS	2.6
1	B	25	PRO	2.5
1	C	366	LEU	2.5
1	C	25	PRO	2.4
1	D	22	GLY	2.3
1	C	211	ASP	2.3
1	B	185	GLY	2.2
1	C	353	ALA	2.1

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	OIC	J	6	11/12	0.94	0.08	25,30,37,39	0
2	ALC	K	3	11/12	0.95	0.10	23,25,30,30	0
2	OIC	I	6	11/12	0.95	0.08	29,31,34,37	0
2	ALC	H	3	11/12	0.95	0.11	28,31,36,37	0
2	ALC	I	3	11/12	0.96	0.10	19,22,28,30	0
2	OIC	H	6	11/12	0.96	0.08	32,33,39,41	0
2	OIC	K	6	11/12	0.96	0.08	28,29,34,36	0
2	ALC	J	3	11/12	0.97	0.10	22,26,36,36	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	PG4	D	401	13/13	0.76	0.21	57,60,77,79	0
3	PG4	A	401	13/13	0.89	0.14	43,47,62,62	0
3	PG4	B	401	13/13	0.91	0.14	36,41,49,52	0
4	PEG	C	401	7/7	0.92	0.21	24,25,25,33	0
4	PEG	B	402	7/7	0.93	0.29	25,25,25,35	0
4	PEG	A	402	7/7	0.93	0.30	25,25,26,32	0

6.5 Other polymers (i)

There are no such residues in this entry.