



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:28 AM UTC

PDB ID : 7CR6 / pdb_00007cr6
Title : Synechocystis Cas1-Cas2/prespacer binary complex
Authors : Yu, Y.; Chen, Q.
Deposited on : 2020-08-12
Resolution : 3.72 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

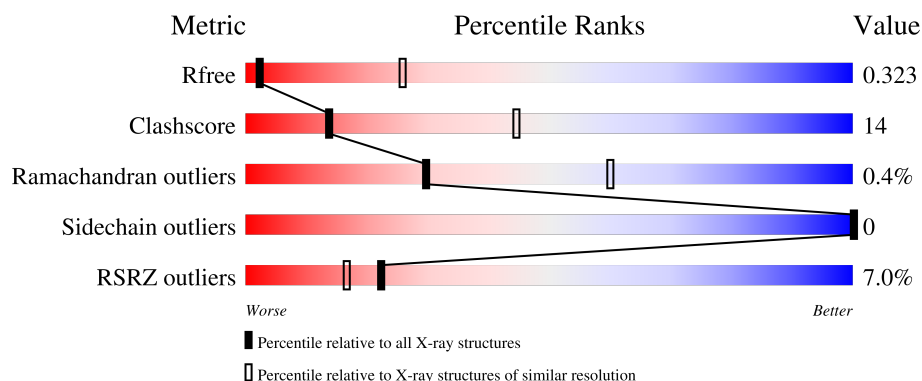
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.72 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	1182 (3.84-3.60)
Clashscore	190562	1222 (3.84-3.60)
Ramachandran outliers	187476	1178 (3.84-3.60)
Sidechain outliers	187428	1174 (3.84-3.60)
RSRZ outliers	180081	1181 (3.84-3.60)

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	336	7% 74% 21% • •
1	B	336	9% 64% 31% • •
1	C	336	5% 65% 31% •
1	D	336	3% 68% 24% 7%
2	E	105	7% 67% 21% 12%

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Mol	Chain	Length	Quality of chain
2	F	105	5% 69% 19% 12%
3	G	36	36% 8% 64% 28%
4	H	36	6% 19% 50% 31%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 24088 atoms, of which 11235 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 CRISPR-associated endonuclease Cas1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	322	Total	C	H	N	O	S	0	0	0
			5072	1661	2484	462	460	5			
1	B	323	Total	C	H	N	O	S	0	0	0
			5038	1663	2445	461	464	5			
1	C	325	Total	C	H	N	O	S	0	0	0
			4966	1674	2354	466	466	6			
1	D	313	Total	C	H	N	O	S	0	0	0
			4970	1622	2445	451	447	5			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	expression tag	UNP Q6ZEI2
A	-9	ALA	-	expression tag	UNP Q6ZEI2
A	-8	SER	-	expression tag	UNP Q6ZEI2
A	-7	GLY	-	expression tag	UNP Q6ZEI2
A	-6	SER	-	expression tag	UNP Q6ZEI2
A	-5	GLY	-	expression tag	UNP Q6ZEI2
A	-4	THR	-	expression tag	UNP Q6ZEI2
A	-3	GLY	-	expression tag	UNP Q6ZEI2
A	-2	SER	-	expression tag	UNP Q6ZEI2
A	-1	GLY	-	expression tag	UNP Q6ZEI2
A	0	SER	-	expression tag	UNP Q6ZEI2
B	-10	GLY	-	expression tag	UNP Q6ZEI2
B	-9	ALA	-	expression tag	UNP Q6ZEI2
B	-8	SER	-	expression tag	UNP Q6ZEI2
B	-7	GLY	-	expression tag	UNP Q6ZEI2
B	-6	SER	-	expression tag	UNP Q6ZEI2
B	-5	GLY	-	expression tag	UNP Q6ZEI2
B	-4	THR	-	expression tag	UNP Q6ZEI2
B	-3	GLY	-	expression tag	UNP Q6ZEI2
B	-2	SER	-	expression tag	UNP Q6ZEI2
B	-1	GLY	-	expression tag	UNP Q6ZEI2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	expression tag	UNP Q6ZEI2
C	-10	GLY	-	expression tag	UNP Q6ZEI2
C	-9	ALA	-	expression tag	UNP Q6ZEI2
C	-8	SER	-	expression tag	UNP Q6ZEI2
C	-7	GLY	-	expression tag	UNP Q6ZEI2
C	-6	SER	-	expression tag	UNP Q6ZEI2
C	-5	GLY	-	expression tag	UNP Q6ZEI2
C	-4	THR	-	expression tag	UNP Q6ZEI2
C	-3	GLY	-	expression tag	UNP Q6ZEI2
C	-2	SER	-	expression tag	UNP Q6ZEI2
C	-1	GLY	-	expression tag	UNP Q6ZEI2
C	0	SER	-	expression tag	UNP Q6ZEI2
D	-10	GLY	-	expression tag	UNP Q6ZEI2
D	-9	ALA	-	expression tag	UNP Q6ZEI2
D	-8	SER	-	expression tag	UNP Q6ZEI2
D	-7	GLY	-	expression tag	UNP Q6ZEI2
D	-6	SER	-	expression tag	UNP Q6ZEI2
D	-5	GLY	-	expression tag	UNP Q6ZEI2
D	-4	THR	-	expression tag	UNP Q6ZEI2
D	-3	GLY	-	expression tag	UNP Q6ZEI2
D	-2	SER	-	expression tag	UNP Q6ZEI2
D	-1	GLY	-	expression tag	UNP Q6ZEI2
D	0	SER	-	expression tag	UNP Q6ZEI2

- Molecule 2 is a protein called CRISPR-associated endoribonuclease Cas2 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	92	Total	C	H	N	O	S	0	0	0
			1494	485	748	123	135	3			
2	F	92	Total	C	H	N	O	S	0	0	0
			1505	485	759	123	135	3			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	GLY	-	expression tag	UNP Q6ZEI1
E	-9	ALA	-	expression tag	UNP Q6ZEI1
E	-8	SER	-	expression tag	UNP Q6ZEI1
E	-7	GLY	-	expression tag	UNP Q6ZEI1
E	-6	SER	-	expression tag	UNP Q6ZEI1
E	-5	GLY	-	expression tag	UNP Q6ZEI1
E	-4	THR	-	expression tag	UNP Q6ZEI1

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-3	GLY	-	expression tag	UNP Q6ZEI1
E	-2	SER	-	expression tag	UNP Q6ZEI1
E	-1	GLY	-	expression tag	UNP Q6ZEI1
E	0	SER	-	expression tag	UNP Q6ZEI1
F	-10	GLY	-	expression tag	UNP Q6ZEI1
F	-9	ALA	-	expression tag	UNP Q6ZEI1
F	-8	SER	-	expression tag	UNP Q6ZEI1
F	-7	GLY	-	expression tag	UNP Q6ZEI1
F	-6	SER	-	expression tag	UNP Q6ZEI1
F	-5	GLY	-	expression tag	UNP Q6ZEI1
F	-4	THR	-	expression tag	UNP Q6ZEI1
F	-3	GLY	-	expression tag	UNP Q6ZEI1
F	-2	SER	-	expression tag	UNP Q6ZEI1
F	-1	GLY	-	expression tag	UNP Q6ZEI1
F	0	SER	-	expression tag	UNP Q6ZEI1

- Molecule 3 is a DNA chain called DNA (36-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	26	Total	C	N	O	P	0	0	0
			526	252	81	167	26			

- Molecule 4 is a DNA chain called DNA (36-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	25	Total	C	N	O	P	0	0	0
			516	243	102	146	25			

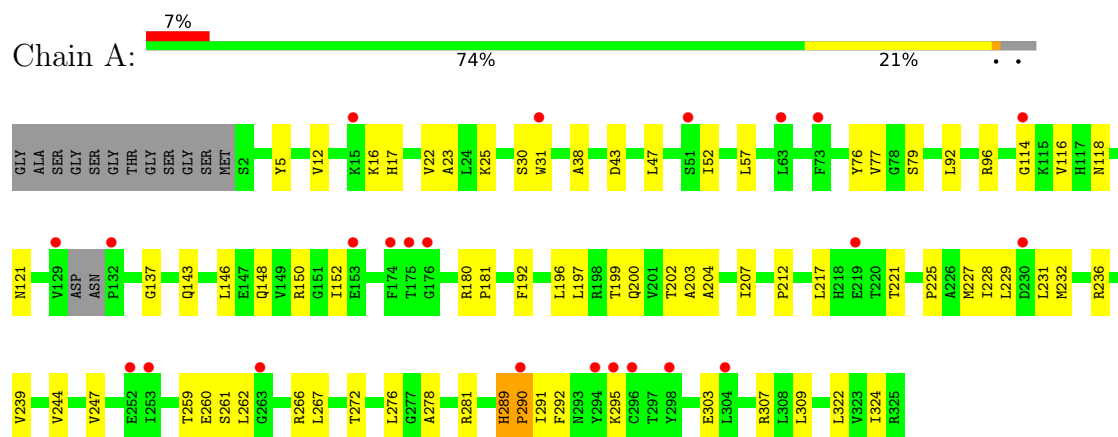
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		

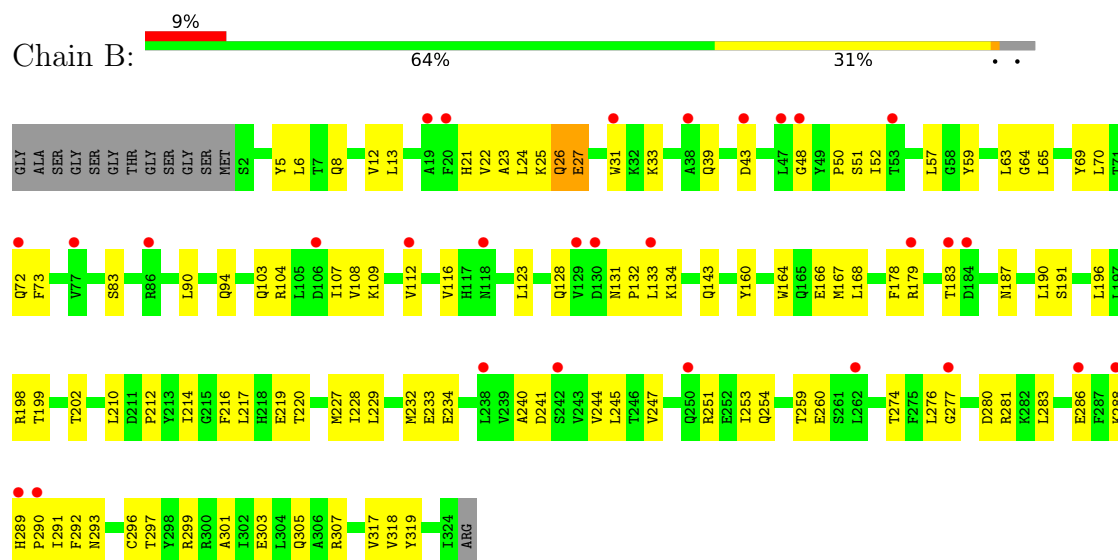
3 Residue-property plots [i](#)

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.

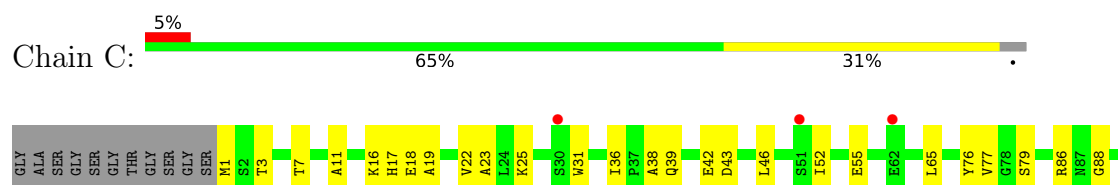
• Molecule 1: CRISPR-associated endonuclease Cas1

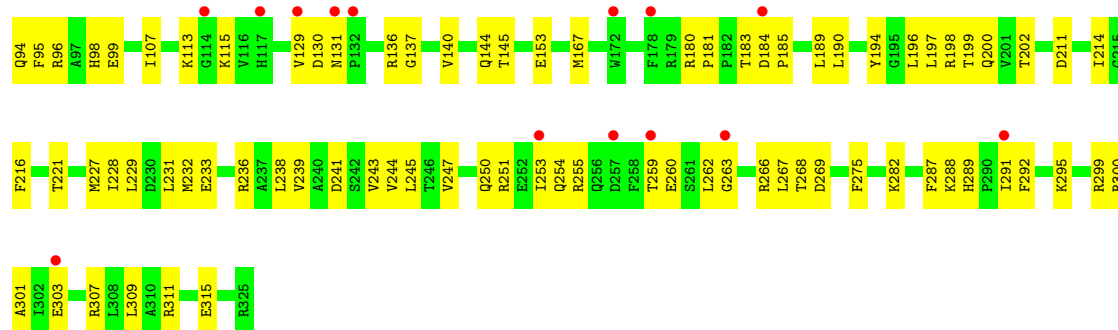


• Molecule 1: CRISPR-associated endonuclease Cas1

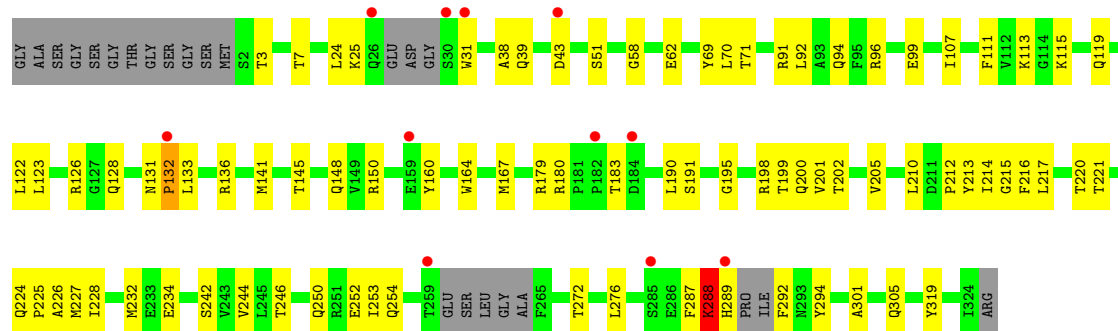


• Molecule 1: CRISPR-associated endonuclease Cas1

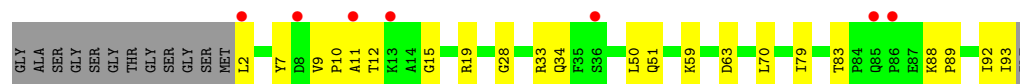




• Molecule 1: CRISPR-associated endonuclease Cas1



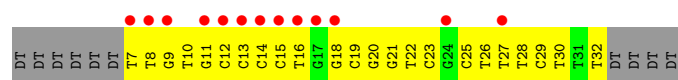
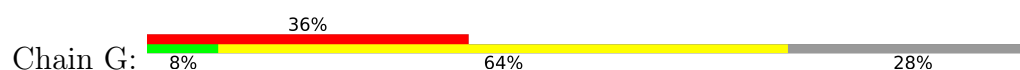
• Molecule 2: CRISPR-associated endonuclease Cas2 1



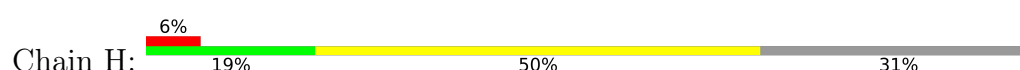
• Molecule 2: CRISPR-associated endonuclease Cas2 1

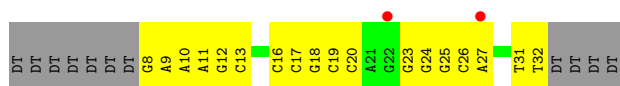


• Molecule 3: DNA (36-MER)



• Molecule 4: DNA (36-MER)





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	125.56Å 215.16Å 191.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.81 – 3.72 33.81 – 3.72	Depositor EDS
% Data completeness (in resolution range)	91.9 (33.81-3.72) 91.8 (33.81-3.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 3.76Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.276 , 0.324 0.277 , 0.323	Depositor DCC
R_{free} test set	1058 reflections (3.85%)	wwPDB-VP
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 25.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.035 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.79	EDS
Total number of atoms	24088	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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](#)

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	0.14	0/2646	0.40	0/3577
1	B	0.17	0/2652	0.43	0/3589
1	C	0.16	0/2671	0.40	0/3613
1	D	0.15	0/2580	0.37	0/3486
2	E	0.12	0/761	0.33	0/1028
2	F	0.16	0/761	0.43	0/1028
3	G	0.16	0/584	0.35	0/899
4	H	0.15	0/580	0.28	0/893
All	All	0.15	0/13235	0.39	0/18113

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
1	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	288	LYS	Peptide

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	2588	2484	2621	54	0
1	B	2593	2445	2618	90	1
1	C	2612	2354	2643	92	1
1	D	2525	2445	2554	77	0
2	E	746	748	767	25	0
2	F	746	759	767	14	0
3	G	526	0	298	29	0
4	H	516	0	279	17	0
5	A	1	0	0	0	0
All	All	12853	11235	12547	367	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (367) 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)
4:H:12:DG:H2''	4:H:13:DC:H5''	1.31	1.07
1:B:25:LYS:HD2	1:B:25:LYS:O	1.63	0.97
1:D:288:LYS:HD2	1:D:289:HIS:H	1.34	0.90
1:C:46:LEU:HD21	1:C:52:ILE:HD11	1.52	0.90
1:B:25:LYS:HE2	1:B:31:TRP:HE1	1.36	0.90
1:C:99:GLU:OE2	1:D:96:ARG:NH1	2.06	0.89
1:D:24:LEU:HD23	1:D:25:LYS:H	1.37	0.88
3:G:25:DC:H2'	3:G:26:DT:H72	1.56	0.87
1:C:288:LYS:HE2	1:C:295:LYS:HE3	1.55	0.87
1:C:43:ASP:OD2	1:C:200:GLN:NE2	2.06	0.87
1:D:150:ARG:NH2	1:D:217:LEU:O	2.09	0.85
1:C:254:GLN:OE1	1:C:255:ARG:N	2.09	0.85
1:D:201:VAL:O	1:D:205:VAL:HG23	1.79	0.82
1:B:303:GLU:OE2	1:B:307:ARG:NH1	2.14	0.81
1:B:25:LYS:O	1:B:26:GLN:HG2	1.80	0.81
1:B:234:GLU:OE2	1:B:319:TYR:OH	1.97	0.80
1:C:253:ILE:HD12	1:C:254:GLN:H	1.46	0.80
1:C:255:ARG:HG2	1:C:255:ARG:O	1.82	0.78
1:A:143:GLN:NE2	1:A:148:GLN:OE1	2.17	0.78
1:D:212:PRO:HA	1:D:227:MET:HB3	1.64	0.78
1:D:123:LEU:HD12	1:D:132:PRO:HB3	1.66	0.76
1:C:253:ILE:HD12	1:C:254:GLN:N	2.02	0.74
1:C:288:LYS:HA	1:C:295:LYS:HA	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:GLU:N	1:A:260:GLU:OE1	2.21	0.73
2:E:33:ARG:NH1	2:F:66:CYS:SG	2.62	0.73
1:C:211:ASP:OD1	1:D:91:ARG:NH1	2.22	0.73
4:H:12:DG:C2'	4:H:13:DC:H5''	2.14	0.73
1:B:216:PHE:CE2	1:B:227:MET:HE1	2.24	0.71
1:C:115:LYS:NZ	1:C:153:GLU:OE1	2.23	0.71
1:D:288:LYS:HD2	1:D:289:HIS:N	2.06	0.70
1:B:241:ASP:O	1:B:245:LEU:HD13	1.90	0.70
1:C:262:LEU:HD13	1:C:262:LEU:O	1.91	0.70
1:B:25:LYS:O	1:B:26:GLN:CG	2.39	0.70
1:B:43:ASP:OD2	1:B:307:ARG:NH2	2.24	0.70
1:C:46:LEU:HD21	1:C:52:ILE:CD1	2.21	0.70
1:A:114:GLY:O	1:A:118:ASN:ND2	2.25	0.70
1:A:303:GLU:OE2	1:A:307:ARG:NH2	2.25	0.70
1:B:253:ILE:HG22	1:B:254:GLN:H	1.55	0.70
1:C:303:GLU:OE2	1:C:307:ARG:NH2	2.24	0.69
1:D:198:ARG:O	1:D:202:THR:OG1	2.09	0.69
3:G:9:DG:H2'	3:G:10:DT:H72	1.74	0.69
1:D:288:LYS:O	1:D:289:HIS:HB2	1.92	0.68
4:H:10:DA:H2'	4:H:11:DA:C8	2.28	0.68
1:C:190:LEU:O	1:C:194:TYR:N	2.27	0.68
1:B:212:PRO:HA	1:B:227:MET:HB3	1.75	0.67
1:C:236:ARG:O	1:C:241:ASP:HB2	1.95	0.67
1:D:24:LEU:CD2	1:D:25:LYS:H	2.07	0.67
2:E:7:TYR:OH	3:G:21:DG:OP1	2.10	0.67
1:B:198:ARG:HA	1:B:232:MET:HE2	1.77	0.67
1:B:51:SER:OG	4:H:9:DA:H5'	1.95	0.66
1:A:76:TYR:OH	1:A:79:SER:OG	2.12	0.66
1:B:228:ILE:HG22	1:B:232:MET:HE3	1.76	0.66
1:B:128:GLN:NE2	1:B:166:GLU:OE1	2.29	0.66
3:G:15:DC:H2''	3:G:16:DT:H71	1.78	0.66
1:D:288:LYS:HG2	1:D:294:TYR:CB	2.26	0.65
1:C:189:LEU:HD22	1:C:244:VAL:HG13	1.77	0.65
1:C:55:GLU:N	1:C:55:GLU:OE1	2.29	0.65
2:E:9:VAL:HG12	2:E:10:PRO:HD2	1.78	0.65
1:B:164:TRP:O	1:B:168:LEU:HD12	1.97	0.65
1:C:260:GLU:HB2	1:C:266:ARG:NH2	2.11	0.64
2:F:29:TYR:CG	2:F:50:LEU:HD12	2.33	0.64
1:C:1:MET:HB3	1:C:42:GLU:OE1	1.98	0.64
1:A:22:VAL:HG11	2:E:93:ILE:CD1	2.28	0.63
1:C:227:MET:HE1	1:C:309:LEU:HD23	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:LEU:HG	1:C:232:MET:HG3	1.81	0.63
1:D:113:LYS:HD2	1:D:141:MET:HE3	1.79	0.63
1:C:190:LEU:CD2	1:C:244:VAL:HG21	2.29	0.62
1:C:287:PHE:CE1	1:C:289:HIS:HA	2.34	0.62
2:E:88:LYS:H	2:E:88:LYS:HD3	1.64	0.62
1:B:259:THR:HG22	1:B:260:GLU:H	1.64	0.62
3:G:12:DC:H2''	3:G:13:DC:O5'	2.00	0.62
1:D:25:LYS:O	1:D:31:TRP:HB2	1.99	0.61
1:C:196:LEU:O	1:C:199:THR:HG22	1.99	0.61
1:C:227:MET:HE1	1:C:309:LEU:CD2	2.30	0.61
3:G:25:DC:H2'	3:G:26:DT:C7	2.30	0.60
1:B:73:PHE:HB2	3:G:30:DT:H2'	1.82	0.60
1:D:288:LYS:HG3	1:D:292:PHE:O	2.02	0.60
1:B:190:LEU:HD12	1:B:244:VAL:HG21	1.83	0.60
2:E:2:LEU:N	2:E:70:LEU:O	2.34	0.60
1:A:25:LYS:HG2	1:A:30:SER:HB2	1.84	0.59
1:B:69:TYR:C	1:B:70:LEU:HD12	2.27	0.59
1:B:288:LYS:HE3	1:B:292:PHE:HB3	1.85	0.59
2:E:88:LYS:HD3	2:E:88:LYS:N	2.17	0.59
1:A:180:ARG:HG3	1:A:181:PRO:HD3	1.85	0.59
1:D:287:PHE:O	1:D:288:LYS:HB2	2.00	0.59
2:E:11:ALA:HB2	3:G:19:DC:H6	1.68	0.59
3:G:18:DG:N2	4:H:20:DC:O2	2.37	0.58
1:C:17:HIS:CE1	3:G:10:DT:H3'	2.39	0.58
1:B:94:GLN:HE21	1:B:227:MET:HE3	1.69	0.57
1:D:160:TYR:O	1:D:164:TRP:NE1	2.38	0.57
1:A:77:VAL:O	1:B:83:SER:N	2.37	0.56
1:B:108:VAL:HG22	1:B:217:LEU:HB2	1.87	0.56
1:A:197:LEU:HD13	1:A:239:VAL:HG21	1.88	0.56
1:B:8:GLN:HE22	1:B:25:LYS:HG2	1.71	0.56
2:E:51:GLN:OE1	2:F:81:TYR:OH	2.23	0.56
1:B:214:ILE:O	1:B:227:MET:HE2	2.06	0.56
1:A:43:ASP:OD2	1:A:200:GLN:NE2	2.38	0.56
1:A:121:ASN:ND2	1:A:324:ILE:O	2.35	0.55
1:B:25:LYS:CE	1:B:31:TRP:HE1	2.14	0.55
1:A:12:VAL:HG12	1:A:23:ALA:HB3	1.87	0.55
1:D:183:THR:O	1:D:183:THR:HG22	2.06	0.55
1:B:191:SER:O	3:G:32:DT:H73	2.07	0.55
2:F:50:LEU:HD23	2:F:50:LEU:C	2.31	0.55
1:D:94:GLN:NE2	1:D:227:MET:HE3	2.22	0.55
1:B:164:TRP:O	1:B:167:MET:N	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:34:GLN:NE2	2:E:83:THR:HG21	2.22	0.54
1:D:224:GLN:O	1:D:226:ALA:N	2.40	0.54
1:C:288:LYS:CE	1:C:295:LYS:HE3	2.34	0.53
1:D:288:LYS:CB	1:D:294:TYR:H	2.20	0.53
1:A:196:LEU:O	1:A:199:THR:HG22	2.08	0.53
1:C:266:ARG:HG3	1:C:267:LEU:O	2.09	0.53
1:C:250:GLN:N	1:C:250:GLN:OE1	2.42	0.53
1:C:259:THR:OG1	1:C:263:GLY:O	2.18	0.53
1:C:38:ALA:O	1:C:65:LEU:HD11	2.08	0.53
1:C:260:GLU:HA	1:C:260:GLU:OE1	2.08	0.53
1:C:287:PHE:HE1	1:C:289:HIS:HA	1.74	0.52
1:C:214:ILE:O	1:C:227:MET:HG2	2.10	0.52
1:C:221:THR:HG22	1:C:221:THR:O	2.09	0.52
4:H:16:DC:H2''	4:H:17:DC:O5'	2.09	0.52
1:C:190:LEU:HD22	1:C:244:VAL:HG21	1.92	0.52
1:D:69:TYR:C	1:D:70:LEU:HD12	2.34	0.52
1:D:220:THR:O	1:D:221:THR:HG23	2.08	0.52
1:D:253:ILE:HG22	1:D:254:GLN:H	1.74	0.52
4:H:18:DG:H2''	4:H:19:DC:O5'	2.09	0.52
1:A:244:VAL:HA	1:A:247:VAL:HG22	1.89	0.52
1:B:6:LEU:HD12	1:B:13:LEU:HD11	1.91	0.52
1:C:76:TYR:OH	1:C:79:SER:OG	2.12	0.52
1:D:234:GLU:OE2	1:D:319:TYR:OH	2.28	0.52
4:H:24:DG:H2''	4:H:25:DG:O5'	2.10	0.52
1:A:192:PHE:CE2	1:A:196:LEU:HD21	2.45	0.52
4:H:17:DC:H2''	4:H:18:DG:C8	2.44	0.52
1:B:178:PHE:N	1:B:187:ASN:OD1	2.39	0.52
1:C:136:ARG:O	1:C:140:VAL:HG23	2.09	0.52
1:C:288:LYS:HE2	1:C:295:LYS:CE	2.32	0.51
1:A:272:THR:O	1:A:276:LEU:HG	2.11	0.51
1:B:296:CYS:SG	1:B:297:THR:N	2.82	0.51
2:F:5:ILE:HD13	2:F:26:LEU:HD21	1.92	0.51
1:A:5:TYR:CG	1:A:276:LEU:HD13	2.45	0.51
1:A:203:ALA:O	1:A:207:ILE:HG13	2.11	0.51
1:B:291:ILE:O	1:B:291:ILE:HG22	2.10	0.51
1:D:7:THR:O	1:D:7:THR:HG22	2.11	0.51
1:D:190:LEU:O	1:D:191:SER:C	2.54	0.51
3:G:13:DC:C2'	3:G:14:DC:H5'	2.41	0.50
1:B:59:TYR:O	1:B:63:LEU:HD23	2.10	0.50
1:B:196:LEU:O	1:B:199:THR:OG1	2.29	0.50
1:C:259:THR:O	1:C:260:GLU:HG2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:289:HIS:CE1	1:C:291:ILE:HD12	2.47	0.50
1:D:69:TYR:O	1:D:70:LEU:HD12	2.10	0.50
1:C:94:GLN:HE21	1:C:227:MET:HG3	1.75	0.50
1:D:225:PRO:HB2	1:D:228:ILE:HD12	1.91	0.50
2:E:92:ILE:O	2:E:93:ILE:HG23	2.11	0.50
1:B:22:VAL:HG12	1:B:24:LEU:HD12	1.93	0.50
1:C:19:ALA:HB2	2:F:28:GLY:CA	2.42	0.50
1:C:229:LEU:HA	1:C:232:MET:HE2	1.93	0.50
1:B:6:LEU:CD1	1:B:13:LEU:HD11	2.41	0.50
1:A:259:THR:O	1:A:259:THR:HG22	2.11	0.50
1:D:51:SER:OG	3:G:8:DT:OP1	2.25	0.50
1:B:90:LEU:HD11	1:B:210:LEU:HD23	1.94	0.50
1:D:133:LEU:HD23	1:D:136:ARG:HG3	1.94	0.50
3:G:19:DC:H2"	3:G:20:DG:O5'	2.11	0.50
2:E:9:VAL:HG12	2:E:10:PRO:CD	2.42	0.49
1:A:289:HIS:CD2	1:A:290:PRO:HD2	2.47	0.49
1:B:22:VAL:CG1	1:B:24:LEU:HD12	2.42	0.49
1:A:150:ARG:NH2	1:A:217:LEU:O	2.45	0.49
1:A:116:VAL:HG21	1:A:137:GLY:HA2	1.94	0.49
1:D:215:GLY:HA2	1:D:227:MET:HE2	1.93	0.49
1:B:103:GLN:O	1:B:107:ILE:HD12	2.12	0.49
1:C:7:THR:HG22	1:C:7:THR:O	2.13	0.49
1:B:39:GLN:O	1:B:39:GLN:NE2	2.46	0.49
1:C:95:PHE:HA	1:C:98:HIS:HB3	1.95	0.49
1:A:52:ILE:HG21	1:A:57:LEU:HD13	1.95	0.49
1:C:289:HIS:HE1	1:C:291:ILE:HD12	1.77	0.49
1:D:128:GLN:OE1	1:D:167:MET:HB2	2.13	0.49
1:B:52:ILE:HD13	1:B:57:LEU:HD11	1.95	0.49
1:B:198:ARG:CA	1:B:232:MET:HE2	2.42	0.48
1:D:253:ILE:HG22	1:D:254:GLN:N	2.27	0.48
1:C:94:GLN:NE2	1:C:227:MET:HG3	2.28	0.48
1:C:43:ASP:OD1	1:C:43:ASP:N	2.43	0.48
1:A:92:LEU:O	1:A:96:ARG:HG3	2.13	0.48
1:D:115:LYS:NZ	1:D:119:GLN:OE1	2.45	0.48
1:B:220:THR:HG23	1:B:220:THR:O	2.13	0.48
1:C:311:ARG:O	1:C:315:GLU:N	2.37	0.48
1:D:221:THR:O	1:D:224:GLN:N	2.45	0.48
2:F:2:LEU:HD13	2:F:75:VAL:HG11	1.95	0.48
3:G:11:DG:H2"	3:G:12:DC:C6	2.49	0.48
1:B:317:VAL:HG12	1:B:318:VAL:N	2.29	0.48
1:D:126:ARG:CB	1:D:167:MET:HE3	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:PRO:HA	1:A:227:MET:HB3	1.95	0.48
1:A:289:HIS:CD2	1:A:322:LEU:HD11	2.48	0.48
1:B:25:LYS:C	1:B:26:GLN:HG2	2.39	0.48
1:B:259:THR:HG22	1:B:260:GLU:N	2.26	0.48
1:C:1:MET:O	1:C:1:MET:HG3	2.12	0.48
1:B:164:TRP:CZ3	1:B:245:LEU:HD11	2.49	0.48
1:B:5:TYR:OH	1:B:303:GLU:OE1	2.19	0.48
1:B:198:ARG:O	1:B:202:THR:OG1	2.20	0.48
1:B:240:ALA:O	1:B:244:VAL:HG23	2.13	0.48
2:F:44:VAL:O	2:F:47:PHE:N	2.46	0.47
1:D:242:SER:O	1:D:246:THR:OG1	2.21	0.47
3:G:11:DG:H4'	3:G:11:DG:OP1	2.14	0.47
1:D:107:ILE:HG22	1:D:111:PHE:HE2	1.79	0.47
1:A:22:VAL:HG11	2:E:93:ILE:HD11	1.96	0.47
1:A:289:HIS:O	1:A:291:ILE:N	2.48	0.47
1:A:292:PHE:HB2	1:A:295:LYS:HD3	1.96	0.47
1:D:216:PHE:CE2	1:D:227:MET:HE1	2.49	0.47
1:D:272:THR:O	1:D:276:LEU:HD13	2.14	0.47
1:D:288:LYS:HG2	1:D:294:TYR:HB2	1.96	0.47
2:E:9:VAL:O	2:E:19:ARG:NH2	2.48	0.47
3:G:7:DT:O4'	3:G:8:DT:H73	2.14	0.47
1:B:183:THR:HG22	1:B:183:THR:O	2.15	0.47
2:F:5:ILE:HD13	2:F:26:LEU:CD2	2.44	0.47
1:A:92:LEU:HD11	1:B:214:ILE:HD12	1.97	0.46
3:G:9:DG:H2'	3:G:10:DT:C7	2.43	0.46
4:H:23:DG:H2''	4:H:24:DG:O5'	2.14	0.46
1:A:92:LEU:CD1	1:B:214:ILE:HD12	2.45	0.46
2:E:59:LYS:N	2:E:63:ASP:OD2	2.43	0.46
1:C:129:VAL:HG13	1:C:130:ASP:H	1.80	0.46
1:C:238:LEU:HD12	1:C:239:VAL:HG13	1.97	0.46
1:D:252:GLU:OE1	1:D:252:GLU:N	2.48	0.46
1:D:288:LYS:CD	1:D:289:HIS:H	2.16	0.46
1:C:107:ILE:HG21	1:C:216:PHE:CD2	2.50	0.46
1:C:268:THR:HG22	1:C:269:ASP:O	2.14	0.46
1:B:48:GLY:C	1:B:50:PRO:HD3	2.40	0.46
1:C:25:LYS:HB3	1:C:31:TRP:HA	1.97	0.46
1:A:197:LEU:HD13	1:A:239:VAL:CG2	2.45	0.46
1:D:71:THR:HG21	4:H:31:DT:OP1	2.16	0.46
1:A:231:LEU:O	1:A:231:LEU:HD23	2.16	0.46
1:B:52:ILE:HD13	1:B:57:LEU:CD1	2.46	0.46
1:D:288:LYS:HG2	1:D:294:TYR:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:22:DT:H2''	3:G:23:DC:C6	2.50	0.46
1:C:289:HIS:CE1	1:C:291:ILE:HB	2.51	0.46
1:B:21:HIS:CE1	1:B:33:LYS:HD3	2.51	0.46
1:C:232:MET:O	1:C:232:MET:HG2	2.14	0.46
1:C:16:LYS:O	1:C:18:GLU:N	2.39	0.45
1:C:202:THR:HG22	1:C:202:THR:O	2.16	0.45
3:G:27:DT:H2''	3:G:28:DT:O5'	2.16	0.45
1:A:146:LEU:HD21	1:A:150:ARG:HH21	1.81	0.45
1:C:167:MET:HE3	1:C:245:LEU:HD21	1.97	0.45
1:B:109:LYS:NZ	1:B:143:GLN:O	2.41	0.45
1:D:38:ALA:O	1:D:39:GLN:HG2	2.17	0.45
1:C:11:ALA:HA	1:C:23:ALA:O	2.16	0.45
2:E:11:ALA:HB2	3:G:19:DC:C6	2.51	0.45
1:B:228:ILE:CG2	1:B:232:MET:HE3	2.45	0.45
1:D:25:LYS:HB3	1:D:31:TRP:HE3	1.81	0.45
1:B:25:LYS:O	1:B:26:GLN:CB	2.64	0.45
1:C:184:ASP:OD2	1:C:184:ASP:C	2.59	0.45
1:C:130:ASP:O	1:C:131:ASN:C	2.59	0.45
1:B:288:LYS:HB3	1:B:293:ASN:O	2.16	0.44
1:D:202:THR:HG23	1:D:228:ILE:HD11	1.99	0.44
1:D:276:LEU:N	1:D:276:LEU:HD12	2.32	0.44
2:E:79:ILE:O	2:F:67:ILE:N	2.49	0.44
2:F:31:LYS:HG3	2:F:32:TRP:H	1.82	0.44
1:A:38:ALA:O	2:E:28:GLY:O	2.36	0.44
1:C:289:HIS:O	1:C:292:PHE:O	2.34	0.44
1:D:228:ILE:O	1:D:232:MET:HG3	2.17	0.44
2:E:88:LYS:O	2:E:88:LYS:HG2	2.17	0.44
1:D:250:GLN:OE1	1:D:250:GLN:N	2.51	0.44
2:E:10:PRO:HA	3:G:19:DC:H5'	1.98	0.44
3:G:8:DT:H2''	3:G:9:DG:C8	2.53	0.44
1:D:3:THR:HG22	1:D:43:ASP:HB2	1.99	0.44
1:C:22:VAL:CG2	1:C:36:ILE:HD12	2.48	0.44
1:A:261:SER:O	1:A:262:LEU:C	2.60	0.44
1:B:276:LEU:N	1:B:276:LEU:HD12	2.33	0.44
1:B:283:LEU:HB3	1:B:299:ARG:HB2	1.99	0.44
1:C:243:VAL:HG23	1:C:282:LYS:HG2	2.00	0.44
1:C:254:GLN:O	1:C:255:ARG:HB3	2.18	0.44
1:B:63:LEU:HG	1:B:65:LEU:HD13	2.00	0.44
1:A:225:PRO:O	1:A:229:LEU:HD13	2.18	0.43
1:B:112:VAL:O	1:B:116:VAL:HG12	2.18	0.43
1:B:123:LEU:HD21	1:B:160:TYR:HE1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:131:ASN:N	1:D:132:PRO:CD	2.81	0.43
2:E:12:THR:HG23	2:E:15:GLY:H	1.83	0.43
1:A:204:ALA:HB1	1:A:303:GLU:HA	2.00	0.43
1:B:247:VAL:HG12	1:B:251:ARG:HD2	2.00	0.43
1:B:253:ILE:HG22	1:B:254:GLN:N	2.27	0.43
1:B:94:GLN:NE2	1:B:227:MET:HE3	2.33	0.43
1:B:164:TRP:HB3	1:B:168:LEU:HD11	2.00	0.43
1:C:260:GLU:HB2	1:C:266:ARG:HH21	1.82	0.43
4:H:32:DT:O2	4:H:32:DT:O4'	2.36	0.43
1:A:31:TRP:O	1:A:31:TRP:CD1	2.72	0.43
1:B:131:ASN:N	1:B:132:PRO:CD	2.81	0.43
1:C:233:GLU:OE2	1:C:236:ARG:NH1	2.51	0.43
1:D:301:ALA:O	1:D:305:GLN:HG2	2.18	0.43
1:B:290:PRO:O	1:B:292:PHE:N	2.52	0.43
1:C:3:THR:HG23	1:C:43:ASP:CG	2.43	0.43
1:C:88:GLY:C	1:D:214:ILE:HD11	2.44	0.43
1:B:290:PRO:C	1:B:292:PHE:H	2.25	0.43
1:C:167:MET:HE3	1:C:245:LEU:CD2	2.49	0.43
1:A:262:LEU:C	1:A:262:LEU:HD13	2.44	0.43
1:C:144:GLN:O	1:C:145:THR:HG23	2.18	0.43
1:D:123:LEU:HD12	1:D:132:PRO:CB	2.43	0.43
1:A:278:ALA:O	1:A:281:ARG:HB2	2.18	0.42
2:F:19:ARG:O	2:F:22:LEU:N	2.52	0.42
3:G:8:DT:OP1	3:G:8:DT:H6	2.02	0.42
4:H:25:DG:H2''	4:H:26:DC:C6	2.54	0.42
4:H:27:DA:H8	4:H:27:DA:O5'	2.02	0.42
2:E:10:PRO:O	2:E:19:ARG:NH2	2.52	0.42
1:A:232:MET:HE3	1:A:236:ARG:NE	2.35	0.42
1:B:104:ARG:NH2	1:B:219:GLU:OE2	2.51	0.42
1:B:160:TYR:O	1:B:164:TRP:CD1	2.72	0.42
1:B:274:THR:O	1:B:277:GLY:N	2.49	0.42
1:C:77:VAL:HG11	1:D:62:GLU:HG2	2.01	0.42
1:C:88:GLY:O	1:D:214:ILE:HD11	2.20	0.42
1:D:145:THR:HG22	1:D:148:GLN:OE1	2.20	0.42
1:D:160:TYR:O	1:D:164:TRP:CD1	2.73	0.42
1:D:190:LEU:HD12	1:D:244:VAL:HG21	2.01	0.42
1:B:25:LYS:O	1:B:25:LYS:CD	2.50	0.42
1:B:280:ASP:OD1	1:B:281:ARG:N	2.53	0.42
2:E:50:LEU:O	2:E:50:LEU:HD23	2.18	0.42
1:B:12:VAL:HG23	1:B:23:ALA:HB3	2.02	0.42
1:D:199:THR:OG1	1:D:200:GLN:OE1	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:21:DG:H1	4:H:16:DC:H42	1.67	0.42
1:C:198:ARG:HG2	1:C:228:ILE:CG2	2.50	0.42
1:D:122:LEU:HD22	1:D:160:TYR:OH	2.20	0.42
3:G:25:DC:H2'	3:G:26:DT:C5	2.54	0.42
1:C:189:LEU:HG	1:C:275:PHE:HE1	1.85	0.42
2:E:88:LYS:N	2:E:89:PRO:HD3	2.35	0.42
1:A:16:LYS:O	1:A:17:HIS:C	2.62	0.42
1:B:297:THR:O	1:B:301:ALA:N	2.47	0.42
1:C:300:ARG:O	1:C:301:ALA:C	2.63	0.42
1:D:195:GLY:O	1:D:198:ARG:HB3	2.20	0.42
1:D:205:VAL:HG13	1:D:210:LEU:HB2	2.02	0.42
2:F:6:ILE:HG22	2:F:37:VAL:HG22	2.01	0.42
2:F:58:ILE:HG23	2:F:63:ASP:OD2	2.19	0.42
1:A:22:VAL:HG11	2:E:93:ILE:HD12	2.00	0.41
1:D:58:GLY:O	1:D:62:GLU:HG3	2.20	0.41
1:D:92:LEU:O	1:D:96:ARG:HG3	2.20	0.41
1:B:290:PRO:C	1:B:292:PHE:N	2.79	0.41
1:C:96:ARG:NE	1:D:99:GLU:OE2	2.53	0.41
1:B:301:ALA:O	1:B:305:GLN:HG2	2.20	0.41
1:C:238:LEU:HD12	1:C:239:VAL:CG1	2.50	0.41
1:D:221:THR:O	1:D:224:GLN:HB2	2.20	0.41
3:G:13:DC:H2''	3:G:14:DC:H5'	2.02	0.41
1:B:289:HIS:HB2	1:B:290:PRO:CD	2.51	0.41
1:A:47:LEU:HD11	1:A:276:LEU:HD11	2.03	0.41
1:B:72:GLN:HB3	3:G:30:DT:H1'	2.03	0.41
1:B:286:GLU:OE2	1:B:288:LYS:HA	2.20	0.41
1:A:5:TYR:CB	1:A:276:LEU:HD13	2.51	0.41
1:A:16:LYS:HG3	1:A:17:HIS:H	1.86	0.41
1:A:202:THR:OG1	1:A:228:ILE:HD13	2.20	0.41
1:A:221:THR:O	1:A:221:THR:HG22	2.20	0.41
1:C:181:PRO:O	1:C:183:THR:N	2.54	0.41
1:D:200:GLN:OE1	1:D:200:GLN:N	2.47	0.41
1:D:288:LYS:HB2	1:D:294:TYR:O	2.21	0.41
1:B:26:GLN:HB3	1:B:27:GLU:H	1.58	0.41
1:C:231:LEU:HD23	1:C:231:LEU:O	2.20	0.41
1:A:148:GLN:O	1:A:152:ILE:HG12	2.21	0.41
1:C:3:THR:OG1	1:C:299:ARG:NH2	2.54	0.41
1:C:180:ARG:HB3	1:C:181:PRO:HD2	2.03	0.41
1:C:184:ASP:OD2	1:C:185:PRO:N	2.53	0.41
1:D:150:ARG:HH21	1:D:217:LEU:HD12	1.86	0.41
4:H:8:DG:H4'	4:H:8:DG:OP2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:ARG:HG3	1:D:213:TYR:CE1	2.56	0.41
1:C:113:LYS:NZ	1:C:137:GLY:O	2.51	0.41
4:H:18:DG:OP2	4:H:18:DG:H8	2.04	0.41
1:A:289:HIS:O	1:A:290:PRO:C	2.63	0.40
1:B:64:GLY:O	1:B:65:LEU:HD12	2.20	0.40
1:B:167:MET:CE	1:B:245:LEU:HD21	2.51	0.40
1:C:247:VAL:O	1:C:251:ARG:O	2.39	0.40
3:G:28:DT:H2'	3:G:29:DC:O4'	2.21	0.40
1:A:227:MET:CE	1:A:309:LEU:HD21	2.52	0.40
1:B:229:LEU:O	1:B:233:GLU:HG2	2.21	0.40
1:A:266:ARG:NE	1:A:267:LEU:O	2.41	0.40
1:D:145:THR:HG22	1:D:148:GLN:CD	2.46	0.40
1:D:179:ARG:O	1:D:180:ARG:C	2.65	0.40
1:B:133:LEU:HD12	1:B:134:LYS:N	2.37	0.40
1:C:18:GLU:O	1:C:39:GLN:NE2	2.50	0.40
1:C:129:VAL:HG13	1:C:130:ASP:N	2.37	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:ARG:O	1:C:144:GLN:HE22[4_555]	1.58	0.02

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	318/336 (95%)	282 (89%)	34 (11%)	2 (1%)	21	52
1	B	321/336 (96%)	282 (88%)	37 (12%)	2 (1%)	21	52
1	C	323/336 (96%)	279 (86%)	44 (14%)	0	100	100
1	D	305/336 (91%)	268 (88%)	35 (12%)	2 (1%)	18	49

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	90/105 (86%)	84 (93%)	6 (7%)	0	100	100
2	F	90/105 (86%)	80 (89%)	10 (11%)	0	100	100
All	All	1447/1554 (93%)	1275 (88%)	166 (12%)	6 (0%)	30	60

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	26	GLN
1	D	288	LYS
1	D	132	PRO
1	A	289	HIS
1	A	290	PRO
1	B	27	GLU

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	272/280 (97%)	272 (100%)	0	100	100
1	B	273/280 (98%)	273 (100%)	0	100	100
1	C	275/280 (98%)	275 (100%)	0	100	100
1	D	266/280 (95%)	266 (100%)	0	100	100
2	E	82/89 (92%)	82 (100%)	0	100	100
2	F	82/89 (92%)	82 (100%)	0	100	100
All	All	1250/1298 (96%)	1250 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	GLN

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Mol	Chain	Res	Type
1	A	148	GLN
1	A	289	HIS
1	A	293	ASN
1	B	8	GLN
1	B	39	GLN
1	B	117	HIS
1	C	289	HIS
1	D	117	HIS
2	F	34	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no 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

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	322/336 (95%)	0.93	23 (7%)	22	16	13, 48, 87, 107	0
1	B	323/336 (96%)	0.85	29 (8%)	15	13	15, 41, 75, 106	0
1	C	325/336 (96%)	0.73	17 (5%)	33	22	9, 33, 75, 135	0
1	D	313/336 (93%)	0.64	11 (3%)	47	30	13, 36, 67, 88	0
2	E	92/105 (87%)	0.83	7 (7%)	20	15	25, 48, 79, 97	0
2	F	92/105 (87%)	0.88	5 (5%)	31	21	13, 47, 63, 69	0
3	G	26/36 (72%)	2.07	13 (50%)	0	1	41, 71, 105, 114	0
4	H	25/36 (69%)	1.53	2 (8%)	18	14	38, 67, 88, 91	0
All	All	1518/1626 (93%)	0.83	107 (7%)	22	17	9, 41, 82, 135	0

All (107) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	30	SER	5.3
1	C	30	SER	4.8
1	C	184	ASP	4.7
1	A	290	PRO	3.9
1	C	253	ILE	3.9
1	B	290	PRO	3.8
1	A	114	GLY	3.7
3	G	15	DC	3.5
1	A	296	CYS	3.5
1	B	289	HIS	3.4
1	A	31	TRP	3.3
1	A	252	GLU	3.3
3	G	16	DT	3.2
1	B	129	VAL	3.1
1	D	26	GLN	3.0
2	E	2	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	178	PHE	2.9
1	B	288	LYS	2.9
4	H	27	DA	2.8
3	G	27	DT	2.8
1	A	295	LYS	2.8
2	F	82	GLY	2.8
1	B	43	ASP	2.8
1	A	298	TYR	2.8
3	G	7	DT	2.7
1	A	219	GLU	2.7
1	A	263	GLY	2.7
1	A	294	TYR	2.7
2	E	11	ALA	2.6
1	C	259	THR	2.6
1	A	253	ILE	2.6
1	D	289	HIS	2.6
3	G	8	DT	2.6
1	B	262	LEU	2.6
1	B	106	ASP	2.6
1	B	130	ASP	2.6
1	C	51	SER	2.6
2	E	13	LYS	2.6
2	E	85	GLN	2.6
3	G	24	DG	2.6
3	G	12	DC	2.5
1	A	175	THR	2.5
1	B	183	THR	2.5
3	G	9	DG	2.5
1	D	182	PRO	2.5
1	A	304	LEU	2.5
1	B	53	THR	2.5
1	C	114	GLY	2.5
1	B	20	PHE	2.4
1	D	31	TRP	2.4
1	B	179	ARG	2.4
1	C	263	GLY	2.4
1	A	51	SER	2.4
3	G	17	DG	2.4
1	A	230	ASP	2.4
1	B	184	ASP	2.4
1	A	176	GLY	2.4
1	A	174	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
3	G	14	DC	2.4
3	G	18	DG	2.4
4	H	22	DG	2.4
1	C	291	ILE	2.4
3	G	13	DC	2.4
1	D	184	ASP	2.4
1	B	31	TRP	2.3
2	F	46	GLN	2.3
1	B	277	GLY	2.3
1	D	259	THR	2.3
1	A	73	PHE	2.3
1	B	242	SER	2.3
1	C	131	ASN	2.3
1	C	62	GLU	2.3
2	E	36	SER	2.3
1	D	43	ASP	2.3
3	G	11	DG	2.3
1	B	72	GLN	2.2
1	C	129	VAL	2.2
1	C	172	TRP	2.2
2	F	7	TYR	2.2
2	F	20	THR	2.2
1	B	38	ALA	2.2
1	B	112	VAL	2.1
2	F	37	VAL	2.1
1	B	286	GLU	2.1
1	A	63	LEU	2.1
1	B	238	LEU	2.1
1	A	132	PRO	2.1
1	D	285	SER	2.1
1	D	132	PRO	2.1
1	B	47	LEU	2.1
1	C	257	ASP	2.1
2	E	8	ASP	2.1
1	B	48	GLY	2.1
1	D	159	GLU	2.1
1	B	19	ALA	2.1
1	B	250	GLN	2.1
1	C	132	PRO	2.1
1	B	86	ARG	2.1
1	A	129	VAL	2.1
2	E	86	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	15	LYS	2.0
1	B	133	LEU	2.0
1	B	118	ASN	2.0
1	A	153	GLU	2.0
1	C	303	GLU	2.0
1	B	77	VAL	2.0
1	C	117	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.