



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:33 PM UTC

PDB ID : 3LYS / pdb_00003lys
Title : Crystal Structure of the N-terminal domain of the Prophage pi2 protein 01 (integrase) from Lactococcus lactis, Northeast Structural Genomics Consortium Target KR124F
Authors : Forouhar, F.; Abashidze, M.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Ciccosanti, C.; Belote, R.L.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2010-02-28
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)

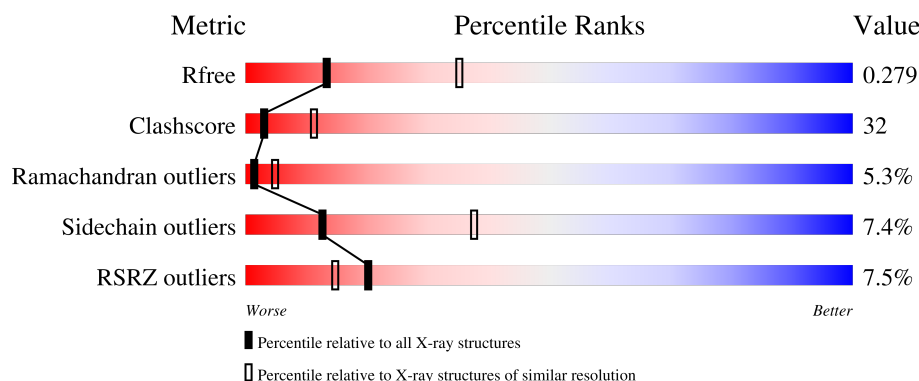
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	112	6% 50% 38% 6% 6%
1	B	112	6% 46% 37% 10% 7%
1	C	112	5% 46% 38% 8% 7%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
1	D	112	8% 42% 39% 13%
1	E	112	7% 46% 41% 5% 7%
1	F	112	8% 44% 43% 7% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5239 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 Prophage pi2 protein 01, integrase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	105	Total	C	N	O	S	Se	0	0	0
			865	553	147	161	1	3			
1	B	104	Total	C	N	O	S	Se	0	0	0
			860	548	148	160	1	3			
1	C	104	Total	C	N	O	S	Se	0	0	0
			855	547	144	160	1	3			
1	D	107	Total	C	N	O	S	Se	0	0	0
			888	566	155	163	1	3			
1	E	104	Total	C	N	O	S	Se	0	0	0
			855	547	144	160	1	3			
1	F	105	Total	C	N	O	S	Se	0	0	0
			865	553	147	161	1	3			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	53	MSE	-	initiating methionine	UNP Q9CGT4
A	137	CYS	PRO	engineered mutation	UNP Q9CGT4
A	157	LEU	-	expression tag	UNP Q9CGT4
A	158	GLU	-	expression tag	UNP Q9CGT4
A	159	HIS	-	expression tag	UNP Q9CGT4
A	160	HIS	-	expression tag	UNP Q9CGT4
A	161	HIS	-	expression tag	UNP Q9CGT4
A	162	HIS	-	expression tag	UNP Q9CGT4
A	163	HIS	-	expression tag	UNP Q9CGT4
A	164	HIS	-	expression tag	UNP Q9CGT4
B	53	MSE	-	initiating methionine	UNP Q9CGT4
B	137	CYS	PRO	engineered mutation	UNP Q9CGT4
B	157	LEU	-	expression tag	UNP Q9CGT4
B	158	GLU	-	expression tag	UNP Q9CGT4
B	159	HIS	-	expression tag	UNP Q9CGT4
B	160	HIS	-	expression tag	UNP Q9CGT4
B	161	HIS	-	expression tag	UNP Q9CGT4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	162	HIS	-	expression tag	UNP Q9CGT4
B	163	HIS	-	expression tag	UNP Q9CGT4
B	164	HIS	-	expression tag	UNP Q9CGT4
C	53	MSE	-	initiating methionine	UNP Q9CGT4
C	137	CYS	PRO	engineered mutation	UNP Q9CGT4
C	157	LEU	-	expression tag	UNP Q9CGT4
C	158	GLU	-	expression tag	UNP Q9CGT4
C	159	HIS	-	expression tag	UNP Q9CGT4
C	160	HIS	-	expression tag	UNP Q9CGT4
C	161	HIS	-	expression tag	UNP Q9CGT4
C	162	HIS	-	expression tag	UNP Q9CGT4
C	163	HIS	-	expression tag	UNP Q9CGT4
C	164	HIS	-	expression tag	UNP Q9CGT4
D	53	MSE	-	initiating methionine	UNP Q9CGT4
D	137	CYS	PRO	engineered mutation	UNP Q9CGT4
D	157	LEU	-	expression tag	UNP Q9CGT4
D	158	GLU	-	expression tag	UNP Q9CGT4
D	159	HIS	-	expression tag	UNP Q9CGT4
D	160	HIS	-	expression tag	UNP Q9CGT4
D	161	HIS	-	expression tag	UNP Q9CGT4
D	162	HIS	-	expression tag	UNP Q9CGT4
D	163	HIS	-	expression tag	UNP Q9CGT4
D	164	HIS	-	expression tag	UNP Q9CGT4
E	53	MSE	-	initiating methionine	UNP Q9CGT4
E	137	CYS	PRO	engineered mutation	UNP Q9CGT4
E	157	LEU	-	expression tag	UNP Q9CGT4
E	158	GLU	-	expression tag	UNP Q9CGT4
E	159	HIS	-	expression tag	UNP Q9CGT4
E	160	HIS	-	expression tag	UNP Q9CGT4
E	161	HIS	-	expression tag	UNP Q9CGT4
E	162	HIS	-	expression tag	UNP Q9CGT4
E	163	HIS	-	expression tag	UNP Q9CGT4
E	164	HIS	-	expression tag	UNP Q9CGT4
F	53	MSE	-	initiating methionine	UNP Q9CGT4
F	137	CYS	PRO	engineered mutation	UNP Q9CGT4
F	157	LEU	-	expression tag	UNP Q9CGT4
F	158	GLU	-	expression tag	UNP Q9CGT4
F	159	HIS	-	expression tag	UNP Q9CGT4
F	160	HIS	-	expression tag	UNP Q9CGT4
F	161	HIS	-	expression tag	UNP Q9CGT4
F	162	HIS	-	expression tag	UNP Q9CGT4
F	163	HIS	-	expression tag	UNP Q9CGT4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	164	HIS	-	expression tag	UNP Q9CGT4

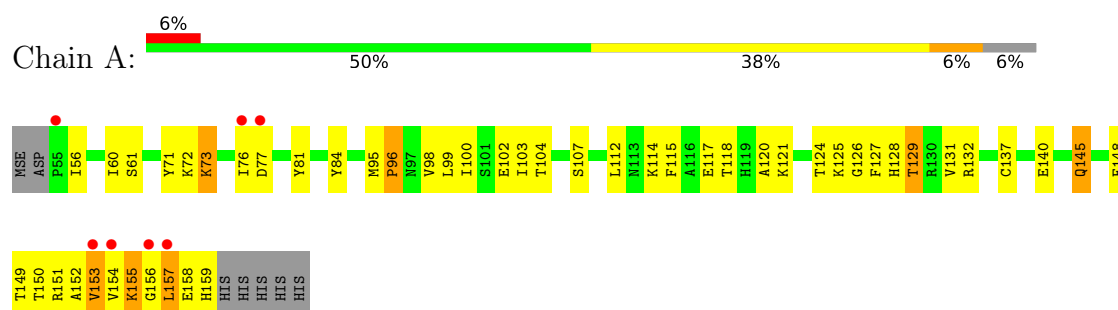
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total O 8 8	0	0
2	B	13	Total O 13 13	0	0
2	C	8	Total O 8 8	0	0
2	D	11	Total O 11 11	0	0
2	E	4	Total O 4 4	0	0
2	F	7	Total O 7 7	0	0

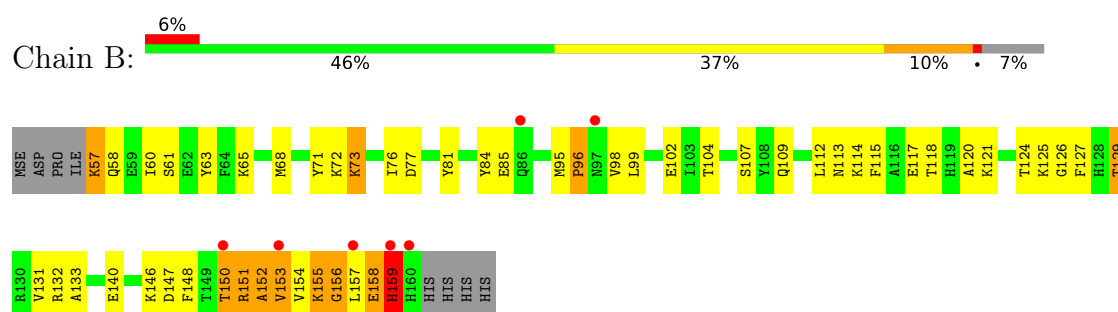
3 Residue-property plots [i](#)

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.

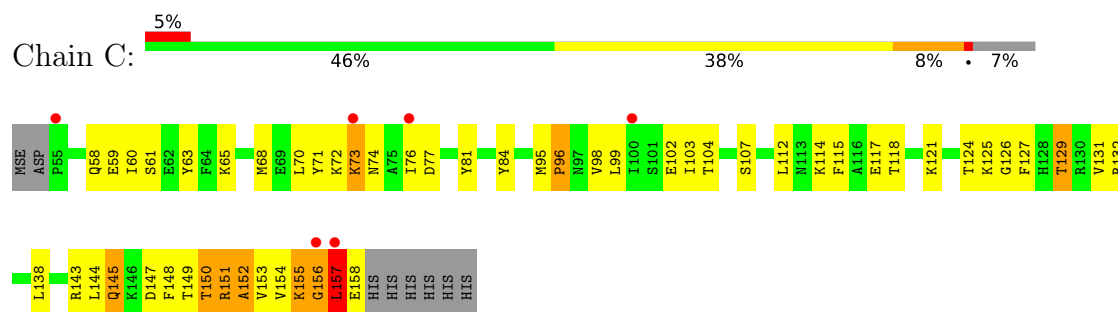
- Molecule 1: Prophage pi2 protein 01, integrase



- Molecule 1: Prophage pi2 protein 01, integrase

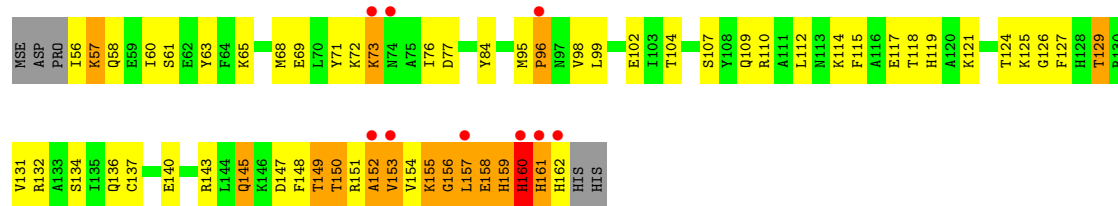


- Molecule 1: Prophage pi2 protein 01, integrase

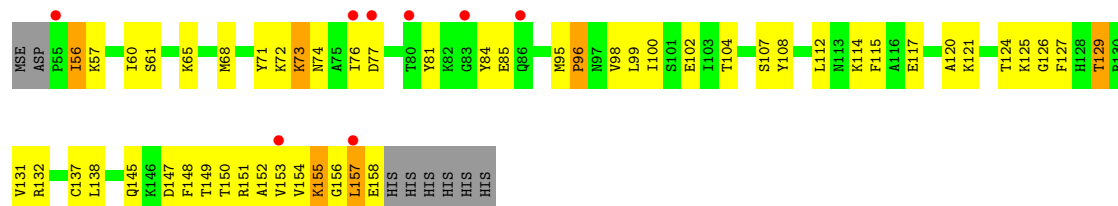


- Molecule 1: Prophage pi2 protein 01, integrase

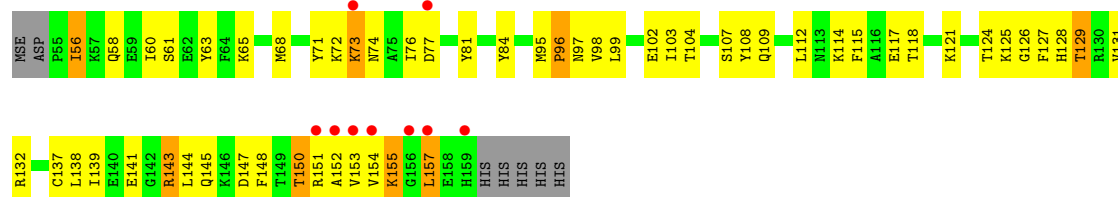
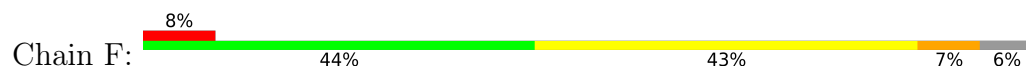




- Molecule 1: Prophage pi2 protein 01, integrase



- Molecule 1: Prophage pi2 protein 01, integrase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.70Å 90.27Å 156.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.80 19.96 – 2.81	Depositor EDS
% Data completeness (in resolution range)	73.2 (19.96-2.80) 84.7 (19.96-2.81)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.08 (at 2.80Å)	Xtriage
Refinement program	CNS 1.2, REFMAC	Depositor
R, R_{free}	0.260 , 0.290 0.253 , 0.279	Depositor DCC
R_{free} test set	1637 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	1.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	5239	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.54	0/879	0.87	1/1174 (0.1%)
1	B	0.57	0/874	0.83	0/1167
1	C	0.56	0/868	0.90	1/1159 (0.1%)
1	D	0.58	0/904	0.98	4/1208 (0.3%)
1	E	0.58	0/868	0.84	0/1159
1	F	0.58	0/879	0.85	1/1174 (0.1%)
All	All	0.57	0/5272	0.88	7/7041 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	110	ARG	NE-CZ-NH2	-11.15	109.17	119.20
1	D	110	ARG	CD-NE-CZ	10.28	138.79	124.40
1	D	110	ARG	NE-CZ-NH1	9.23	130.73	121.50
1	D	160	HIS	N-CA-C	5.71	117.26	107.28
1	F	143	ARG	N-CA-C	-5.48	107.24	114.31
1	A	145	GLN	N-CA-C	5.19	116.79	111.03
1	C	157	LEU	N-CA-C	5.08	114.75	108.19

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	865	0	873	54	0
1	B	860	0	861	63	0
1	C	855	0	866	54	0
1	D	888	0	886	69	0
1	E	855	0	866	50	0
1	F	865	0	873	62	0
2	A	8	0	0	0	0
2	B	13	0	0	4	0
2	C	8	0	0	0	0
2	D	11	0	0	4	0
2	E	4	0	0	0	0
2	F	7	0	0	1	0
All	All	5239	0	5225	329	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 32.

All (329) 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:B:150:THR:HG23	1:B:151:ARG:H	1.26	0.99
1:D:125:LYS:O	1:D:129:THR:HG22	1.62	0.99
1:D:150:THR:HG23	1:D:151:ARG:H	1.28	0.98
1:C:125:LYS:O	1:C:129:THR:HG22	1.64	0.97
1:E:125:LYS:O	1:E:129:THR:HG22	1.64	0.97
1:B:125:LYS:O	1:B:129:THR:HG22	1.65	0.96
1:A:125:LYS:O	1:A:129:THR:HG22	1.66	0.96
1:D:121:LYS:H	1:D:159:HIS:CE1	1.83	0.95
1:A:120:ALA:HA	1:A:158:GLU:HB3	1.50	0.93
1:F:125:LYS:O	1:F:129:THR:HG22	1.69	0.92
1:D:121:LYS:H	1:D:159:HIS:HE1	1.15	0.89
1:A:154:VAL:HG21	1:B:71:TYR:HB3	1.59	0.84
1:F:150:THR:HG23	1:F:151:ARG:H	1.43	0.81
1:C:151:ARG:HG3	1:D:140:GLU:HG3	1.64	0.80
1:A:124:THR:HG21	1:A:156:GLY:O	1.82	0.79
1:D:99:LEU:HB2	1:D:102:GLU:OE2	1.83	0.78
1:A:157:LEU:HD13	1:A:157:LEU:H	1.49	0.77
1:A:150:THR:HG23	1:A:151:ARG:H	1.50	0.76
1:D:161:HIS:HB2	1:E:74:ASN:HB3	1.68	0.76
1:A:140:GLU:HG3	1:F:151:ARG:CD	2.16	0.75
1:F:99:LEU:HB2	1:F:102:GLU:OE2	1.86	0.75
1:C:147:ASP:O	1:C:150:THR:HG22	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ARG:C	1:B:153:VAL:H	1.94	0.74
1:E:99:LEU:HB2	1:E:102:GLU:OE2	1.87	0.73
1:C:99:LEU:HB2	1:C:102:GLU:OE2	1.90	0.72
1:A:99:LEU:HB2	1:A:102:GLU:OE2	1.90	0.72
1:A:150:THR:HG23	1:A:151:ARG:N	2.04	0.71
1:B:155:LYS:H	1:B:155:LYS:HE3	1.55	0.71
1:F:150:THR:HG23	1:F:151:ARG:N	2.05	0.70
1:B:99:LEU:HB2	1:B:102:GLU:OE2	1.91	0.70
1:D:56:ILE:HG22	1:D:143:ARG:O	1.91	0.70
1:D:124:THR:HG21	1:D:156:GLY:O	1.92	0.70
1:A:56:ILE:HG13	1:A:100:ILE:HG21	1.75	0.69
1:C:151:ARG:C	1:C:153:VAL:H	2.00	0.68
1:D:72:LYS:HE2	1:D:84:TYR:OH	1.94	0.68
1:A:132:ARG:HH21	1:A:132:ARG:HG3	1.59	0.68
1:C:132:ARG:HH21	1:C:132:ARG:HG3	1.58	0.68
1:E:132:ARG:HH21	1:E:132:ARG:HG3	1.59	0.68
1:E:151:ARG:C	1:E:153:VAL:H	2.03	0.66
1:F:132:ARG:HH21	1:F:132:ARG:HG3	1.59	0.66
1:C:72:LYS:HE2	1:C:84:TYR:OH	1.95	0.66
1:D:150:THR:HG23	1:D:151:ARG:N	2.05	0.65
1:B:132:ARG:HH21	1:B:132:ARG:HG3	1.59	0.65
1:F:95:MSE:HE3	1:F:98:VAL:HG11	1.79	0.65
1:A:72:LYS:HE2	1:A:84:TYR:OH	1.97	0.65
1:B:152:ALA:C	1:B:154:VAL:H	2.03	0.64
1:B:120:ALA:HA	1:B:158:GLU:HB3	1.78	0.64
1:B:158:GLU:O	1:B:159:HIS:HB2	1.96	0.64
1:A:140:GLU:HG3	1:F:151:ARG:HD3	1.77	0.64
1:D:132:ARG:HH21	1:D:132:ARG:HG3	1.62	0.64
1:B:72:LYS:HE2	1:B:84:TYR:OH	1.97	0.64
1:D:69:GLU:OE2	2:D:13:HOH:O	2.16	0.63
1:D:121:LYS:N	1:D:159:HIS:HE1	1.91	0.63
1:C:138:LEU:HD12	1:C:143:ARG:HD3	1.80	0.62
1:C:95:MSE:HE3	1:C:98:VAL:HG11	1.82	0.62
1:C:148:PHE:CD1	1:C:149:THR:HG23	2.34	0.62
1:B:157:LEU:HB2	1:B:158:GLU:OE2	1.99	0.62
1:B:95:MSE:HE3	1:B:98:VAL:HG11	1.81	0.61
1:D:151:ARG:C	1:D:153:VAL:H	2.07	0.61
1:E:72:LYS:HE2	1:E:84:TYR:OH	1.99	0.61
1:F:72:LYS:HE2	1:F:84:TYR:OH	2.00	0.61
1:D:161:HIS:CB	1:E:74:ASN:HB3	2.29	0.61
1:D:152:ALA:C	1:D:154:VAL:H	2.09	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ARG:C	1:B:153:VAL:N	2.59	0.60
1:C:150:THR:HG23	1:C:151:ARG:H	1.65	0.60
1:A:148:PHE:CE1	1:A:149:THR:HG23	2.36	0.60
1:A:56:ILE:HG23	1:A:56:ILE:O	2.02	0.60
1:F:121:LYS:O	1:F:124:THR:HG22	2.02	0.59
1:E:121:LYS:O	1:E:124:THR:HG22	2.02	0.59
1:A:95:MSE:HE3	1:A:98:VAL:HG11	1.85	0.59
1:F:56:ILE:HG23	1:F:56:ILE:O	2.03	0.59
1:A:148:PHE:CD1	1:A:149:THR:HG23	2.38	0.59
1:C:121:LYS:O	1:C:124:THR:HG22	2.03	0.59
1:B:151:ARG:O	1:B:153:VAL:N	2.35	0.59
1:A:151:ARG:C	1:A:153:VAL:H	2.09	0.58
1:A:151:ARG:HD2	1:B:140:GLU:HG3	1.85	0.58
1:F:151:ARG:HB3	1:F:151:ARG:NH1	2.18	0.58
1:D:145:GLN:OE1	1:D:145:GLN:HA	2.02	0.58
1:B:121:LYS:O	1:B:124:THR:HG22	2.04	0.57
1:A:158:GLU:O	1:A:159:HIS:HB2	2.04	0.57
1:F:132:ARG:NH2	1:F:147:ASP:OD2	2.33	0.57
1:D:57:LYS:HD3	1:D:57:LYS:N	2.19	0.57
1:E:95:MSE:HE3	1:E:98:VAL:HG11	1.87	0.57
1:F:151:ARG:C	1:F:153:VAL:H	2.13	0.57
1:C:157:LEU:HD23	1:C:158:GLU:OE1	2.04	0.57
1:B:114:LYS:O	1:B:117:GLU:HG2	2.04	0.57
1:D:95:MSE:HE3	1:D:98:VAL:HG11	1.87	0.57
1:A:150:THR:CG2	1:A:151:ARG:H	2.18	0.57
1:B:150:THR:CG2	1:B:151:ARG:H	2.09	0.57
1:D:60:ILE:HG21	1:D:95:MSE:HE2	1.86	0.56
1:B:60:ILE:HG21	1:B:95:MSE:HE2	1.86	0.56
1:B:150:THR:HG23	1:B:151:ARG:N	2.06	0.56
1:F:155:LYS:H	1:F:155:LYS:HD2	1.71	0.56
1:F:141:GLU:OE1	1:F:143:ARG:HD2	2.05	0.56
1:D:114:LYS:HE3	2:D:43:HOH:O	2.06	0.56
1:D:121:LYS:O	1:D:124:THR:HG22	2.05	0.55
1:B:121:LYS:NZ	1:B:155:LYS:NZ	2.55	0.55
1:E:150:THR:HG23	1:E:151:ARG:H	1.72	0.55
1:E:124:THR:HG21	1:E:156:GLY:O	2.05	0.55
1:C:60:ILE:HG21	1:C:95:MSE:HE2	1.87	0.55
1:C:58:GLN:OE1	1:C:63:TYR:HB2	2.07	0.55
1:C:114:LYS:O	1:C:117:GLU:HG2	2.08	0.54
1:D:152:ALA:O	1:D:154:VAL:N	2.39	0.54
1:D:157:LEU:HG	1:D:160:HIS:ND1	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:ALA:C	1:B:154:VAL:N	2.65	0.54
1:D:76:ILE:HG13	1:D:77:ASP:N	2.22	0.54
1:E:60:ILE:HG23	1:E:61:SER:N	2.22	0.54
1:F:60:ILE:HG21	1:F:95:MSE:HE2	1.89	0.54
1:F:108:TYR:OH	1:F:128:HIS:ND1	2.29	0.54
1:C:115:PHE:CE1	1:C:124:THR:HA	2.42	0.54
1:E:155:LYS:HD2	1:E:155:LYS:O	2.06	0.54
1:B:60:ILE:HG23	1:B:61:SER:N	2.23	0.54
1:B:126:GLY:O	1:B:129:THR:HG23	2.08	0.53
1:F:95:MSE:CE	1:F:98:VAL:HG11	2.38	0.53
1:F:76:ILE:HG13	1:F:77:ASP:N	2.23	0.53
1:C:60:ILE:HG23	1:C:61:SER:N	2.22	0.53
1:A:155:LYS:H	1:A:155:LYS:HD2	1.73	0.53
1:A:76:ILE:HG13	1:A:77:ASP:N	2.23	0.53
1:B:76:ILE:HG13	1:B:77:ASP:N	2.23	0.53
1:D:159:HIS:N	1:D:159:HIS:ND1	2.55	0.53
1:B:104:THR:H	1:B:107:SER:HB2	1.72	0.53
1:A:121:LYS:O	1:A:124:THR:HG22	2.09	0.53
1:C:104:THR:H	1:C:107:SER:HB2	1.74	0.53
1:E:76:ILE:HG13	1:E:77:ASP:N	2.24	0.53
1:F:132:ARG:HG3	1:F:132:ARG:NH2	2.24	0.52
1:A:114:LYS:O	1:A:117:GLU:HG2	2.09	0.52
1:E:108:TYR:CE1	1:E:149:THR:HG22	2.45	0.52
1:A:132:ARG:HG3	1:A:132:ARG:NH2	2.25	0.52
1:B:132:ARG:HG3	1:B:132:ARG:NH2	2.24	0.52
1:A:60:ILE:HG21	1:A:95:MSE:HE2	1.90	0.52
1:D:115:PHE:CE1	1:D:124:THR:HA	2.44	0.52
1:E:132:ARG:HG3	1:E:132:ARG:NH2	2.25	0.52
1:A:95:MSE:CE	1:A:98:VAL:HG11	2.39	0.52
1:D:114:LYS:O	1:D:117:GLU:HG2	2.10	0.52
1:C:95:MSE:CE	1:C:98:VAL:HG11	2.40	0.52
1:E:158:GLU:HA	1:E:158:GLU:OE1	2.10	0.52
1:F:115:PHE:CE1	1:F:124:THR:HA	2.45	0.52
1:A:104:THR:H	1:A:107:SER:HB2	1.74	0.52
1:E:114:LYS:O	1:E:117:GLU:HG2	2.10	0.52
1:C:76:ILE:HG13	1:C:77:ASP:N	2.25	0.52
1:E:154:VAL:HG21	1:F:71:TYR:HB3	1.90	0.52
1:F:114:LYS:O	1:F:117:GLU:HG2	2.10	0.52
1:A:60:ILE:HG23	1:A:61:SER:N	2.26	0.51
1:B:115:PHE:CE1	1:B:124:THR:HA	2.46	0.51
1:D:60:ILE:HG23	1:D:61:SER:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:147:ASP:O	1:E:150:THR:HG22	2.10	0.51
1:F:104:THR:H	1:F:107:SER:HB2	1.75	0.51
1:D:57:LYS:O	1:D:58:GLN:HB2	2.10	0.51
1:E:56:ILE:HG13	1:E:100:ILE:HG21	1.92	0.51
1:C:132:ARG:HG3	1:C:132:ARG:NH2	2.25	0.51
1:E:60:ILE:HG21	1:E:95:MSE:HE2	1.92	0.51
1:D:104:THR:H	1:D:107:SER:HB2	1.76	0.51
2:B:20:HOH:O	1:C:70:LEU:HD23	2.09	0.51
1:C:127:PHE:O	1:C:131:VAL:HG23	2.11	0.51
1:D:148:PHE:CE1	1:D:149:THR:HG22	2.45	0.51
1:C:156:GLY:C	1:C:157:LEU:HD13	2.36	0.51
1:E:145:GLN:OE1	1:E:145:GLN:HA	2.10	0.51
1:A:140:GLU:HG3	1:F:151:ARG:HD2	1.91	0.51
1:B:152:ALA:O	1:B:154:VAL:N	2.44	0.51
1:C:103:ILE:HB	1:C:148:PHE:CE2	2.46	0.51
1:E:95:MSE:CE	1:E:98:VAL:HG11	2.41	0.51
1:F:60:ILE:HG23	1:F:61:SER:N	2.26	0.50
1:B:155:LYS:HE3	1:B:155:LYS:N	2.25	0.50
1:D:126:GLY:O	1:D:129:THR:HG23	2.11	0.50
1:C:58:GLN:HG2	1:C:59:GLU:N	2.27	0.50
1:E:104:THR:H	1:E:107:SER:HB2	1.77	0.50
1:D:119:HIS:O	1:D:158:GLU:HB2	2.12	0.50
1:B:147:ASP:O	1:B:150:THR:HG22	2.12	0.50
1:D:132:ARG:HG3	1:D:132:ARG:NH2	2.26	0.50
1:F:151:ARG:HB3	1:F:151:ARG:HH11	1.74	0.50
1:F:157:LEU:H	1:F:157:LEU:HD22	1.76	0.50
1:A:145:GLN:HA	1:A:145:GLN:OE1	2.11	0.49
1:D:155:LYS:HD2	1:D:155:LYS:O	2.13	0.49
1:A:115:PHE:CE1	1:A:124:THR:HA	2.48	0.49
1:B:155:LYS:HD2	1:B:155:LYS:O	2.12	0.49
1:E:115:PHE:CE1	1:E:124:THR:HA	2.47	0.49
1:C:121:LYS:NZ	1:C:155:LYS:NZ	2.61	0.49
1:C:126:GLY:O	1:C:129:THR:HG23	2.13	0.49
1:C:144:LEU:HD12	1:C:145:GLN:H	1.77	0.49
1:C:151:ARG:C	1:C:153:VAL:N	2.70	0.48
1:E:126:GLY:O	1:E:129:THR:HG23	2.12	0.48
1:B:58:GLN:OE1	1:B:63:TYR:HB2	2.13	0.48
1:B:146:LYS:NZ	2:B:23:HOH:O	2.46	0.48
1:B:148:PHE:C	1:B:150:THR:H	2.20	0.48
1:C:121:LYS:HZ1	1:C:155:LYS:HZ1	1.62	0.48
1:E:151:ARG:C	1:E:153:VAL:N	2.70	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:126:GLY:O	1:F:129:THR:HG23	2.13	0.48
1:D:127:PHE:O	1:D:131:VAL:HG23	2.14	0.48
1:E:72:LYS:O	1:E:73:LYS:C	2.56	0.47
1:D:109:GLN:NE2	1:D:154:VAL:HB	2.29	0.47
1:D:158:GLU:HA	1:D:161:HIS:NE2	2.30	0.47
1:B:95:MSE:CE	1:B:98:VAL:HG11	2.43	0.47
1:D:76:ILE:HG13	1:D:77:ASP:H	1.79	0.47
1:E:76:ILE:HG13	1:E:77:ASP:H	1.80	0.47
1:C:151:ARG:NH2	1:D:136:GLN:HB2	2.29	0.47
1:B:76:ILE:HG13	1:B:77:ASP:H	1.80	0.47
1:D:96:PRO:HD2	2:D:10:HOH:O	2.13	0.47
1:A:126:GLY:O	1:A:129:THR:HG23	2.15	0.47
1:E:127:PHE:O	1:E:131:VAL:HG23	2.15	0.47
1:F:72:LYS:O	1:F:73:LYS:C	2.57	0.47
1:B:72:LYS:O	1:B:73:LYS:C	2.57	0.47
1:C:154:VAL:HG21	1:D:71:TYR:HB3	1.96	0.47
1:B:155:LYS:O	1:B:155:LYS:CE	2.63	0.46
1:E:153:VAL:O	1:E:155:LYS:HG3	2.15	0.46
1:F:157:LEU:N	1:F:157:LEU:HD13	2.29	0.46
1:D:157:LEU:HB2	1:D:158:GLU:H	1.58	0.46
1:E:73:LYS:HB3	1:E:74:ASN:H	1.60	0.46
1:F:139:ILE:HD13	1:F:144:LEU:HB3	1.97	0.46
1:A:72:LYS:O	1:A:73:LYS:C	2.58	0.46
1:D:95:MSE:CE	1:D:98:VAL:HG11	2.45	0.46
1:C:151:ARG:HH22	1:D:136:GLN:HB2	1.81	0.46
1:A:76:ILE:HG13	1:A:77:ASP:H	1.79	0.46
1:E:121:LYS:HA	1:E:124:THR:HG22	1.97	0.46
1:A:71:TYR:HB3	1:F:154:VAL:HG21	1.98	0.46
1:B:158:GLU:O	1:B:159:HIS:CB	2.61	0.46
1:F:127:PHE:O	1:F:131:VAL:HG23	2.16	0.46
1:B:151:ARG:HE	1:B:151:ARG:HB3	1.52	0.46
1:C:72:LYS:O	1:C:73:LYS:C	2.59	0.45
1:D:72:LYS:HE2	1:D:84:TYR:CZ	2.51	0.45
1:A:151:ARG:C	1:A:153:VAL:N	2.74	0.45
1:B:113:ASN:ND2	2:B:20:HOH:O	2.48	0.45
1:B:121:LYS:HZ1	1:B:155:LYS:NZ	2.14	0.45
1:D:151:ARG:C	1:D:153:VAL:N	2.73	0.45
1:F:128:HIS:NE2	1:F:153:VAL:HG23	2.32	0.45
1:F:121:LYS:NZ	1:F:157:LEU:HD12	2.32	0.45
1:F:109:GLN:NE2	1:F:154:VAL:HB	2.32	0.45
1:A:127:PHE:O	1:A:131:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:LYS:O	1:D:73:LYS:C	2.59	0.45
1:A:121:LYS:HD2	1:A:157:LEU:HB3	1.99	0.45
1:D:154:VAL:HG21	1:E:71:TYR:HB3	1.97	0.45
1:D:157:LEU:HD13	1:D:157:LEU:N	2.32	0.45
1:F:76:ILE:HG13	1:F:77:ASP:H	1.80	0.45
1:D:56:ILE:HD12	1:D:57:LYS:N	2.32	0.45
1:F:103:ILE:HG21	1:F:148:PHE:CE1	2.53	0.44
1:E:65:LYS:O	1:E:68:MSE:HG2	2.18	0.44
1:F:58:GLN:OE1	1:F:63:TYR:HB2	2.18	0.44
1:F:73:LYS:HB3	1:F:74:ASN:H	1.59	0.44
1:B:127:PHE:O	1:B:131:VAL:HG23	2.18	0.44
1:C:72:LYS:HE2	1:C:84:TYR:CZ	2.53	0.44
1:C:76:ILE:HG13	1:C:77:ASP:H	1.83	0.44
1:D:161:HIS:HB2	1:E:74:ASN:CB	2.43	0.44
1:F:151:ARG:O	1:F:153:VAL:N	2.50	0.44
1:A:72:LYS:HE2	1:A:84:TYR:CZ	2.53	0.44
1:C:152:ALA:C	1:C:154:VAL:H	2.24	0.44
1:D:152:ALA:C	1:D:154:VAL:N	2.74	0.44
1:B:73:LYS:HA	1:B:81:TYR:CE1	2.53	0.44
1:D:121:LYS:HA	1:D:124:THR:HG22	1.99	0.44
1:F:96:PRO:C	1:F:98:VAL:H	2.26	0.44
1:A:71:TYR:N	1:A:71:TYR:CD1	2.86	0.44
1:E:151:ARG:O	1:E:153:VAL:N	2.51	0.44
1:C:151:ARG:HG3	1:D:140:GLU:CG	2.41	0.43
1:F:148:PHE:C	1:F:150:THR:H	2.26	0.43
1:B:155:LYS:HE3	1:B:155:LYS:O	2.18	0.43
1:A:73:LYS:HA	1:A:81:TYR:CE1	2.54	0.43
1:C:73:LYS:HB3	1:C:74:ASN:H	1.58	0.43
1:B:72:LYS:HE2	1:B:84:TYR:CZ	2.53	0.43
1:B:121:LYS:NZ	1:B:155:LYS:HZ2	2.16	0.43
1:F:56:ILE:HD13	1:F:138:LEU:HD11	2.01	0.43
1:A:121:LYS:HZ2	1:A:155:LYS:HD3	1.84	0.43
1:B:65:LYS:O	1:B:68:MSE:HG2	2.19	0.43
1:C:71:TYR:CD1	1:C:71:TYR:N	2.87	0.43
1:C:154:VAL:HG11	1:D:71:TYR:O	2.19	0.43
1:E:71:TYR:CD1	1:E:71:TYR:N	2.86	0.43
1:F:65:LYS:O	1:F:68:MSE:HG2	2.19	0.43
1:B:71:TYR:CD1	1:B:71:TYR:N	2.87	0.42
1:B:96:PRO:C	1:B:98:VAL:H	2.27	0.42
1:E:96:PRO:C	1:E:98:VAL:H	2.27	0.42
1:F:138:LEU:HD12	1:F:143:ARG:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:LYS:O	1:D:68:MSE:HG2	2.19	0.42
1:D:147:ASP:OD1	1:D:149:THR:HG23	2.20	0.42
1:A:154:VAL:HG11	1:B:71:TYR:O	2.20	0.42
1:B:57:LYS:N	1:B:57:LYS:CD	2.82	0.42
1:E:60:ILE:HG23	1:E:61:SER:H	1.83	0.42
1:F:77:ASP:HB2	2:F:51:HOH:O	2.18	0.42
1:F:151:ARG:HH11	1:F:151:ARG:CB	2.32	0.42
1:A:152:ALA:C	1:A:154:VAL:N	2.78	0.42
1:C:151:ARG:HD3	1:D:137:CYS:HA	2.02	0.42
1:E:138:LEU:HD12	1:E:138:LEU:HA	1.94	0.42
1:F:121:LYS:HZ2	1:F:157:LEU:HD12	1.84	0.42
1:A:120:ALA:HA	1:A:158:GLU:CB	2.37	0.42
1:C:138:LEU:HD12	1:C:138:LEU:HA	1.94	0.42
1:D:71:TYR:CD1	1:D:71:TYR:N	2.87	0.42
1:E:73:LYS:HA	1:E:81:TYR:CE1	2.55	0.42
1:B:155:LYS:H	1:B:155:LYS:CE	2.28	0.42
1:A:96:PRO:C	1:A:98:VAL:H	2.28	0.42
1:A:137:CYS:SG	1:F:151:ARG:NH1	2.93	0.42
1:A:152:ALA:O	1:A:154:VAL:N	2.53	0.42
1:D:96:PRO:C	1:D:98:VAL:H	2.28	0.42
1:C:60:ILE:HG23	1:C:61:SER:H	1.84	0.42
1:C:148:PHE:CE1	1:C:149:THR:HG23	2.55	0.42
1:F:145:GLN:HA	1:F:145:GLN:OE1	2.19	0.42
1:B:81:TYR:O	1:B:85:GLU:HG3	2.19	0.41
1:F:73:LYS:HA	1:F:81:TYR:CE1	2.55	0.41
1:D:157:LEU:HD21	1:D:161:HIS:HB3	2.02	0.41
1:C:157:LEU:HD23	1:C:158:GLU:CD	2.45	0.41
1:F:151:ARG:C	1:F:153:VAL:N	2.79	0.41
1:B:133:ALA:HA	2:B:33:HOH:O	2.20	0.41
1:B:154:VAL:HG11	1:C:71:TYR:O	2.20	0.41
1:E:151:ARG:HD3	1:F:137:CYS:HA	2.01	0.41
1:B:60:ILE:HG23	1:B:61:SER:H	1.85	0.41
1:E:121:LYS:HB2	1:E:157:LEU:CB	2.50	0.41
1:A:128:HIS:O	1:A:129:THR:C	2.63	0.41
1:B:109:GLN:NE2	1:B:152:ALA:O	2.53	0.41
1:C:151:ARG:HE	1:C:151:ARG:HB3	1.56	0.41
1:D:153:VAL:HG23	2:D:15:HOH:O	2.21	0.41
1:A:157:LEU:H	1:A:157:LEU:CD1	2.27	0.41
1:B:155:LYS:O	1:B:156:GLY:O	2.38	0.41
1:C:65:LYS:O	1:C:68:MSE:HG2	2.21	0.41
1:D:58:GLN:OE1	1:D:63:TYR:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:148:PHE:CD1	1:E:149:THR:HG23	2.56	0.41
1:F:71:TYR:CD1	1:F:71:TYR:N	2.89	0.41
1:F:138:LEU:HD12	1:F:138:LEU:HA	1.97	0.41
1:B:121:LYS:HA	1:B:124:THR:HG22	2.02	0.41
1:F:68:MSE:HE2	1:F:81:TYR:CZ	2.56	0.41
1:F:121:LYS:HA	1:F:124:THR:HG22	2.03	0.41
1:E:81:TYR:O	1:E:85:GLU:HG3	2.22	0.40
1:C:96:PRO:C	1:C:98:VAL:H	2.28	0.40
1:D:57:LYS:HD3	1:D:57:LYS:H	1.87	0.40
1:D:131:VAL:O	1:D:134:SER:OG	2.32	0.40
1:E:72:LYS:HE2	1:E:84:TYR:CZ	2.56	0.40
1:F:72:LYS:HE2	1:F:84:TYR:CZ	2.55	0.40
1:F:126:GLY:HA2	1:F:129:THR:CG2	2.52	0.40
1:A:121:LYS:HA	1:A:124:THR:HG22	2.03	0.40
1:C:157:LEU:HD13	1:C:157:LEU:N	2.36	0.40
1:D:151:ARG:NH1	1:E:137:CYS:HB2	2.36	0.40
1:E:120:ALA:HA	1:E:158:GLU:HB2	2.03	0.40
1:A:60:ILE:HB	1:A:103:ILE:HD11	2.03	0.40
1:B:132:ARG:NH2	1:B:132:ARG:CG	2.85	0.40
1:C:73:LYS:HA	1:C:81:TYR:CE1	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	103/112 (92%)	86 (84%)	14 (14%)	3 (3%)	3	13
1	B	102/112 (91%)	85 (83%)	10 (10%)	7 (7%)	1	2
1	C	102/112 (91%)	88 (86%)	9 (9%)	5 (5%)	1	6
1	D	105/112 (94%)	84 (80%)	13 (12%)	8 (8%)	1	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	102/112 (91%)	84 (82%)	14 (14%)	4 (4%)	2	8
1	F	103/112 (92%)	88 (85%)	9 (9%)	6 (6%)	1	4
All	All	617/672 (92%)	515 (84%)	69 (11%)	33 (5%)	1	5

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	LYS
1	A	96	PRO
1	B	73	LYS
1	B	96	PRO
1	B	156	GLY
1	C	73	LYS
1	C	96	PRO
1	C	150	THR
1	D	73	LYS
1	D	96	PRO
1	D	150	THR
1	D	153	VAL
1	E	56	ILE
1	E	73	LYS
1	E	96	PRO
1	F	73	LYS
1	F	96	PRO
1	F	150	THR
1	B	152	ALA
1	B	153	VAL
1	B	159	HIS
1	C	156	GLY
1	B	150	THR
1	C	152	ALA
1	D	161	HIS
1	E	152	ALA
1	F	152	ALA
1	A	153	VAL
1	D	152	ALA
1	D	158	GLU
1	F	97	ASN
1	D	156	GLY
1	F	56	ILE

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	93/96 (97%)	88 (95%)	5 (5%)	20	51
1	B	92/96 (96%)	84 (91%)	8 (9%)	9	30
1	C	92/96 (96%)	85 (92%)	7 (8%)	12	36
1	D	95/96 (99%)	84 (88%)	11 (12%)	5	18
1	E	92/96 (96%)	87 (95%)	5 (5%)	20	51
1	F	93/96 (97%)	88 (95%)	5 (5%)	20	51
All	All	557/576 (97%)	516 (93%)	41 (7%)	13	37

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LEU
1	A	118	THR
1	A	129	THR
1	A	155	LYS
1	A	157	LEU
1	B	57	LYS
1	B	112	LEU
1	B	118	THR
1	B	129	THR
1	B	151	ARG
1	B	155	LYS
1	B	158	GLU
1	B	159	HIS
1	C	112	LEU
1	C	118	THR
1	C	129	THR
1	C	145	GLN
1	C	151	ARG
1	C	155	LYS
1	C	157	LEU
1	D	57	LYS
1	D	112	LEU

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Mol	Chain	Res	Type
1	D	118	THR
1	D	129	THR
1	D	145	GLN
1	D	149	THR
1	D	155	LYS
1	D	157	LEU
1	D	159	HIS
1	D	160	HIS
1	D	162	HIS
1	E	57	LYS
1	E	112	LEU
1	E	129	THR
1	E	155	LYS
1	E	157	LEU
1	F	112	LEU
1	F	118	THR
1	F	129	THR
1	F	155	LYS
1	F	157	LEU

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	58	GLN
1	A	109	GLN
1	B	109	GLN
1	B	113	ASN
1	C	109	GLN
1	D	109	GLN
1	D	159	HIS
1	E	109	GLN
1	F	97	ASN
1	F	109	GLN

5.3.3 RNA ⓘ

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	102/112 (91%)	0.46	7 (6%)	23 16	20, 40, 71, 85	0
1	B	101/112 (90%)	0.53	7 (6%)	23 16	25, 46, 67, 85	0
1	C	101/112 (90%)	0.53	6 (5%)	28 21	26, 49, 77, 90	0
1	D	104/112 (92%)	0.55	9 (8%)	16 11	18, 43, 76, 102	0
1	E	101/112 (90%)	0.55	8 (7%)	18 13	15, 43, 72, 83	0
1	F	102/112 (91%)	0.56	9 (8%)	15 11	17, 44, 76, 103	0
All	All	611/672 (90%)	0.53	46 (7%)	20 15	15, 44, 76, 103	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	157	LEU	6.4
1	D	157	LEU	5.6
1	B	153	VAL	4.3
1	E	153	VAL	4.1
1	C	157	LEU	3.5
1	F	153	VAL	3.4
1	F	157	LEU	3.4
1	E	157	LEU	3.4
1	D	161	HIS	3.3
1	B	86	GLN	3.2
1	A	55	PRO	3.2
1	E	76	ILE	3.2
1	D	153	VAL	3.1
1	C	156	GLY	2.9
1	A	153	VAL	2.9
1	E	77	ASP	2.8
1	A	77	ASP	2.8
1	D	73	LYS	2.7
1	F	159	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	156	GLY	2.5
1	D	162	HIS	2.5
1	D	96	PRO	2.5
1	B	160	HIS	2.5
1	D	160	HIS	2.5
1	C	100	ILE	2.5
1	A	76	ILE	2.4
1	A	154	VAL	2.4
1	F	152	ALA	2.4
1	E	80	THR	2.4
1	B	157	LEU	2.4
1	B	97	ASN	2.3
1	C	76	ILE	2.3
1	B	150	THR	2.3
1	C	55	PRO	2.3
1	E	83	GLY	2.3
1	C	73	LYS	2.3
1	D	152	ALA	2.2
1	F	151	ARG	2.2
1	F	154	VAL	2.2
1	D	74	ASN	2.2
1	F	156	GLY	2.1
1	E	55	PRO	2.1
1	B	159	HIS	2.0
1	F	77	ASP	2.0
1	E	86	GLN	2.0
1	F	73	LYS	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.