



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:55 AM UTC

PDB ID : 4L3F / pdb_00004l3f
Title : Crystal structure of Internalin K (InlK) from *Listeria monocytogenes*
Authors : Neves, D.
Deposited on : 2013-06-05
Resolution : 2.39 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

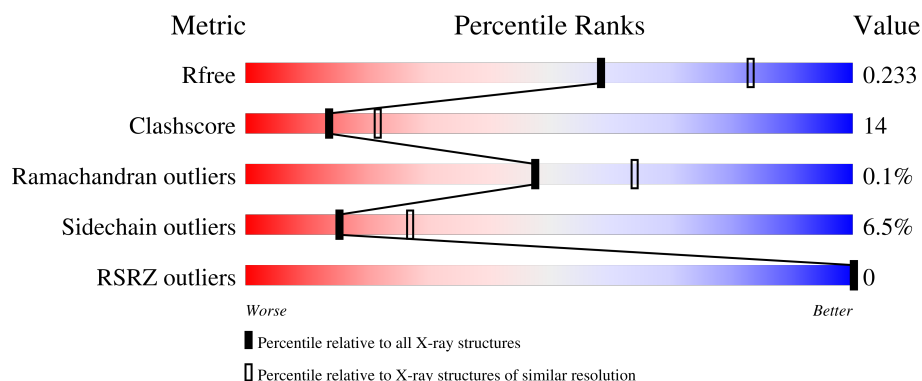
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	347	
1	B	347	
1	C	347	
1	D	347	
1	E	347	

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Mol	Chain	Length	Quality of chain
1	F	347	 64%24%•10%
1	G	347	 62%24%•10%
1	H	347	 59%27%••10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20171 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 Internalin K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	B	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	C	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	D	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	E	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	F	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	G	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			
1	H	311	Total	C	N	O	S	0	0	0
			2436	1528	396	504	8			

There are 176 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	expression tag	UNP Q8Y7I7
A	24	GLY	-	expression tag	UNP Q8Y7I7
A	25	SER	-	expression tag	UNP Q8Y7I7
A	26	SER	-	expression tag	UNP Q8Y7I7
A	27	HIS	-	expression tag	UNP Q8Y7I7
A	28	HIS	-	expression tag	UNP Q8Y7I7
A	29	HIS	-	expression tag	UNP Q8Y7I7
A	30	HIS	-	expression tag	UNP Q8Y7I7
A	31	HIS	-	expression tag	UNP Q8Y7I7
A	32	HIS	-	expression tag	UNP Q8Y7I7
A	33	SER	-	expression tag	UNP Q8Y7I7
A	34	SER	-	expression tag	UNP Q8Y7I7
A	35	GLY	-	expression tag	UNP Q8Y7I7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	36	LEU	-	expression tag	UNP Q8Y7I7
A	37	VAL	-	expression tag	UNP Q8Y7I7
A	38	PRO	-	expression tag	UNP Q8Y7I7
A	39	ARG	-	expression tag	UNP Q8Y7I7
A	40	GLY	-	expression tag	UNP Q8Y7I7
A	41	SER	-	expression tag	UNP Q8Y7I7
A	42	HIS	-	expression tag	UNP Q8Y7I7
A	43	MET	-	expression tag	UNP Q8Y7I7
A	44	ALA	-	expression tag	UNP Q8Y7I7
B	23	MET	-	expression tag	UNP Q8Y7I7
B	24	GLY	-	expression tag	UNP Q8Y7I7
B	25	SER	-	expression tag	UNP Q8Y7I7
B	26	SER	-	expression tag	UNP Q8Y7I7
B	27	HIS	-	expression tag	UNP Q8Y7I7
B	28	HIS	-	expression tag	UNP Q8Y7I7
B	29	HIS	-	expression tag	UNP Q8Y7I7
B	30	HIS	-	expression tag	UNP Q8Y7I7
B	31	HIS	-	expression tag	UNP Q8Y7I7
B	32	HIS	-	expression tag	UNP Q8Y7I7
B	33	SER	-	expression tag	UNP Q8Y7I7
B	34	SER	-	expression tag	UNP Q8Y7I7
B	35	GLY	-	expression tag	UNP Q8Y7I7
B	36	LEU	-	expression tag	UNP Q8Y7I7
B	37	VAL	-	expression tag	UNP Q8Y7I7
B	38	PRO	-	expression tag	UNP Q8Y7I7
B	39	ARG	-	expression tag	UNP Q8Y7I7
B	40	GLY	-	expression tag	UNP Q8Y7I7
B	41	SER	-	expression tag	UNP Q8Y7I7
B	42	HIS	-	expression tag	UNP Q8Y7I7
B	43	MET	-	expression tag	UNP Q8Y7I7
B	44	ALA	-	expression tag	UNP Q8Y7I7
C	23	MET	-	expression tag	UNP Q8Y7I7
C	24	GLY	-	expression tag	UNP Q8Y7I7
C	25	SER	-	expression tag	UNP Q8Y7I7
C	26	SER	-	expression tag	UNP Q8Y7I7
C	27	HIS	-	expression tag	UNP Q8Y7I7
C	28	HIS	-	expression tag	UNP Q8Y7I7
C	29	HIS	-	expression tag	UNP Q8Y7I7
C	30	HIS	-	expression tag	UNP Q8Y7I7
C	31	HIS	-	expression tag	UNP Q8Y7I7
C	32	HIS	-	expression tag	UNP Q8Y7I7
C	33	SER	-	expression tag	UNP Q8Y7I7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	34	SER	-	expression tag	UNP Q8Y7I7
C	35	GLY	-	expression tag	UNP Q8Y7I7
C	36	LEU	-	expression tag	UNP Q8Y7I7
C	37	VAL	-	expression tag	UNP Q8Y7I7
C	38	PRO	-	expression tag	UNP Q8Y7I7
C	39	ARG	-	expression tag	UNP Q8Y7I7
C	40	GLY	-	expression tag	UNP Q8Y7I7
C	41	SER	-	expression tag	UNP Q8Y7I7
C	42	HIS	-	expression tag	UNP Q8Y7I7
C	43	MET	-	expression tag	UNP Q8Y7I7
C	44	ALA	-	expression tag	UNP Q8Y7I7
D	23	MET	-	expression tag	UNP Q8Y7I7
D	24	GLY	-	expression tag	UNP Q8Y7I7
D	25	SER	-	expression tag	UNP Q8Y7I7
D	26	SER	-	expression tag	UNP Q8Y7I7
D	27	HIS	-	expression tag	UNP Q8Y7I7
D	28	HIS	-	expression tag	UNP Q8Y7I7
D	29	HIS	-	expression tag	UNP Q8Y7I7
D	30	HIS	-	expression tag	UNP Q8Y7I7
D	31	HIS	-	expression tag	UNP Q8Y7I7
D	32	HIS	-	expression tag	UNP Q8Y7I7
D	33	SER	-	expression tag	UNP Q8Y7I7
D	34	SER	-	expression tag	UNP Q8Y7I7
D	35	GLY	-	expression tag	UNP Q8Y7I7
D	36	LEU	-	expression tag	UNP Q8Y7I7
D	37	VAL	-	expression tag	UNP Q8Y7I7
D	38	PRO	-	expression tag	UNP Q8Y7I7
D	39	ARG	-	expression tag	UNP Q8Y7I7
D	40	GLY	-	expression tag	UNP Q8Y7I7
D	41	SER	-	expression tag	UNP Q8Y7I7
D	42	HIS	-	expression tag	UNP Q8Y7I7
D	43	MET	-	expression tag	UNP Q8Y7I7
D	44	ALA	-	expression tag	UNP Q8Y7I7
E	23	MET	-	expression tag	UNP Q8Y7I7
E	24	GLY	-	expression tag	UNP Q8Y7I7
E	25	SER	-	expression tag	UNP Q8Y7I7
E	26	SER	-	expression tag	UNP Q8Y7I7
E	27	HIS	-	expression tag	UNP Q8Y7I7
E	28	HIS	-	expression tag	UNP Q8Y7I7
E	29	HIS	-	expression tag	UNP Q8Y7I7
E	30	HIS	-	expression tag	UNP Q8Y7I7
E	31	HIS	-	expression tag	UNP Q8Y7I7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	32	HIS	-	expression tag	UNP Q8Y7I7
E	33	SER	-	expression tag	UNP Q8Y7I7
E	34	SER	-	expression tag	UNP Q8Y7I7
E	35	GLY	-	expression tag	UNP Q8Y7I7
E	36	LEU	-	expression tag	UNP Q8Y7I7
E	37	VAL	-	expression tag	UNP Q8Y7I7
E	38	PRO	-	expression tag	UNP Q8Y7I7
E	39	ARG	-	expression tag	UNP Q8Y7I7
E	40	GLY	-	expression tag	UNP Q8Y7I7
E	41	SER	-	expression tag	UNP Q8Y7I7
E	42	HIS	-	expression tag	UNP Q8Y7I7
E	43	MET	-	expression tag	UNP Q8Y7I7
E	44	ALA	-	expression tag	UNP Q8Y7I7
F	23	MET	-	expression tag	UNP Q8Y7I7
F	24	GLY	-	expression tag	UNP Q8Y7I7
F	25	SER	-	expression tag	UNP Q8Y7I7
F	26	SER	-	expression tag	UNP Q8Y7I7
F	27	HIS	-	expression tag	UNP Q8Y7I7
F	28	HIS	-	expression tag	UNP Q8Y7I7
F	29	HIS	-	expression tag	UNP Q8Y7I7
F	30	HIS	-	expression tag	UNP Q8Y7I7
F	31	HIS	-	expression tag	UNP Q8Y7I7
F	32	HIS	-	expression tag	UNP Q8Y7I7
F	33	SER	-	expression tag	UNP Q8Y7I7
F	34	SER	-	expression tag	UNP Q8Y7I7
F	35	GLY	-	expression tag	UNP Q8Y7I7
F	36	LEU	-	expression tag	UNP Q8Y7I7
F	37	VAL	-	expression tag	UNP Q8Y7I7
F	38	PRO	-	expression tag	UNP Q8Y7I7
F	39	ARG	-	expression tag	UNP Q8Y7I7
F	40	GLY	-	expression tag	UNP Q8Y7I7
F	41	SER	-	expression tag	UNP Q8Y7I7
F	42	HIS	-	expression tag	UNP Q8Y7I7
F	43	MET	-	expression tag	UNP Q8Y7I7
F	44	ALA	-	expression tag	UNP Q8Y7I7
G	23	MET	-	expression tag	UNP Q8Y7I7
G	24	GLY	-	expression tag	UNP Q8Y7I7
G	25	SER	-	expression tag	UNP Q8Y7I7
G	26	SER	-	expression tag	UNP Q8Y7I7
G	27	HIS	-	expression tag	UNP Q8Y7I7
G	28	HIS	-	expression tag	UNP Q8Y7I7
G	29	HIS	-	expression tag	UNP Q8Y7I7

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Chain	Residue	Modelled	Actual	Comment	Reference
G	30	HIS	-	expression tag	UNP Q8Y7I7
G	31	HIS	-	expression tag	UNP Q8Y7I7
G	32	HIS	-	expression tag	UNP Q8Y7I7
G	33	SER	-	expression tag	UNP Q8Y7I7
G	34	SER	-	expression tag	UNP Q8Y7I7
G	35	GLY	-	expression tag	UNP Q8Y7I7
G	36	LEU	-	expression tag	UNP Q8Y7I7
G	37	VAL	-	expression tag	UNP Q8Y7I7
G	38	PRO	-	expression tag	UNP Q8Y7I7
G	39	ARG	-	expression tag	UNP Q8Y7I7
G	40	GLY	-	expression tag	UNP Q8Y7I7
G	41	SER	-	expression tag	UNP Q8Y7I7
G	42	HIS	-	expression tag	UNP Q8Y7I7
G	43	MET	-	expression tag	UNP Q8Y7I7
G	44	ALA	-	expression tag	UNP Q8Y7I7
H	23	MET	-	expression tag	UNP Q8Y7I7
H	24	GLY	-	expression tag	UNP Q8Y7I7
H	25	SER	-	expression tag	UNP Q8Y7I7
H	26	SER	-	expression tag	UNP Q8Y7I7
H	27	HIS	-	expression tag	UNP Q8Y7I7
H	28	HIS	-	expression tag	UNP Q8Y7I7
H	29	HIS	-	expression tag	UNP Q8Y7I7
H	30	HIS	-	expression tag	UNP Q8Y7I7
H	31	HIS	-	expression tag	UNP Q8Y7I7
H	32	HIS	-	expression tag	UNP Q8Y7I7
H	33	SER	-	expression tag	UNP Q8Y7I7
H	34	SER	-	expression tag	UNP Q8Y7I7
H	35	GLY	-	expression tag	UNP Q8Y7I7
H	36	LEU	-	expression tag	UNP Q8Y7I7
H	37	VAL	-	expression tag	UNP Q8Y7I7
H	38	PRO	-	expression tag	UNP Q8Y7I7
H	39	ARG	-	expression tag	UNP Q8Y7I7
H	40	GLY	-	expression tag	UNP Q8Y7I7
H	41	SER	-	expression tag	UNP Q8Y7I7
H	42	HIS	-	expression tag	UNP Q8Y7I7
H	43	MET	-	expression tag	UNP Q8Y7I7
H	44	ALA	-	expression tag	UNP Q8Y7I7

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ni 1 1	0	0
2	F	1	Total Ni 1 1	0	0

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	1	Total Na 1 1	0	0
3	H	1	Total Na 1 1	0	0

- Molecule 4 is water.

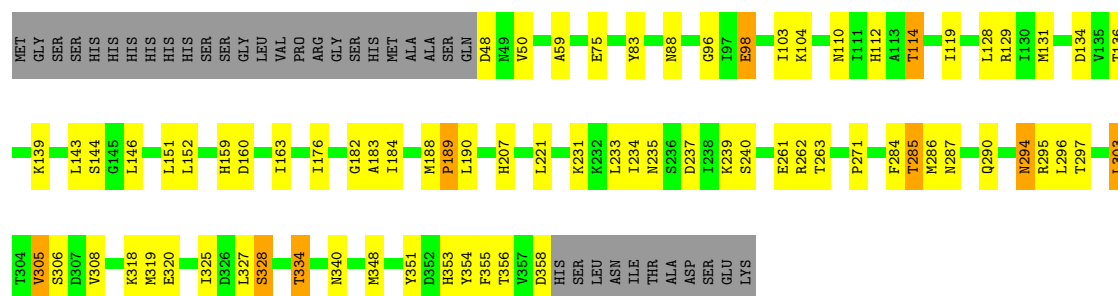
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	96	Total O 96 96	0	0
4	B	93	Total O 93 93	0	0
4	C	99	Total O 99 99	0	0
4	D	95	Total O 95 95	0	0
4	E	59	Total O 59 59	0	0
4	F	87	Total O 87 87	0	0
4	G	65	Total O 65 65	0	0
4	H	85	Total O 85 85	0	0

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.

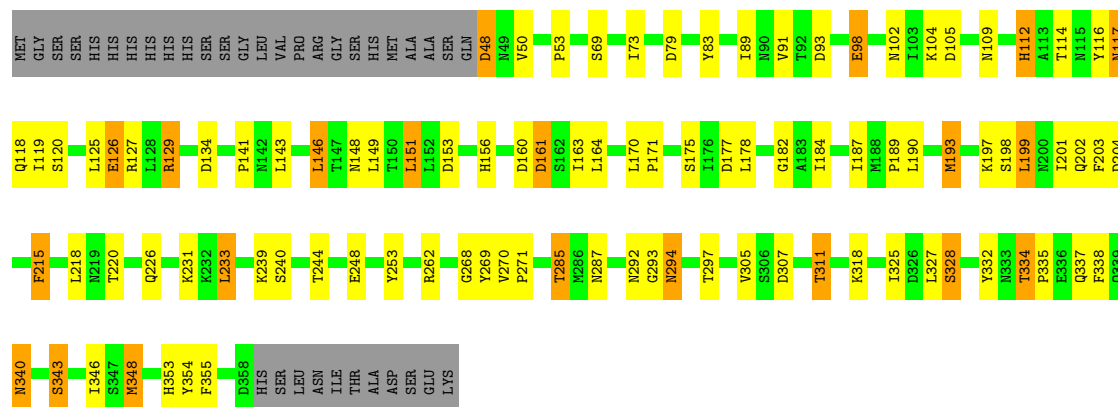
• Molecule 1: Internalin K

Chain A: 



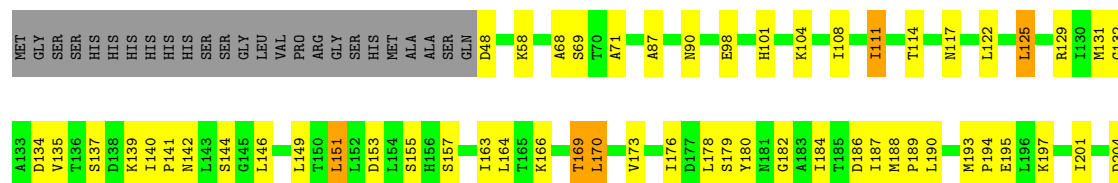
• Molecule 1: Internalin K

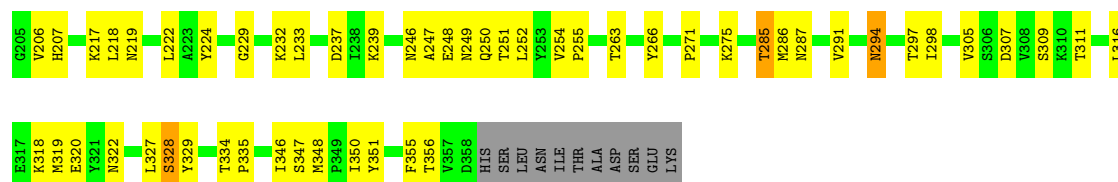
Chain B: 



• Molecule 1: Internalin K

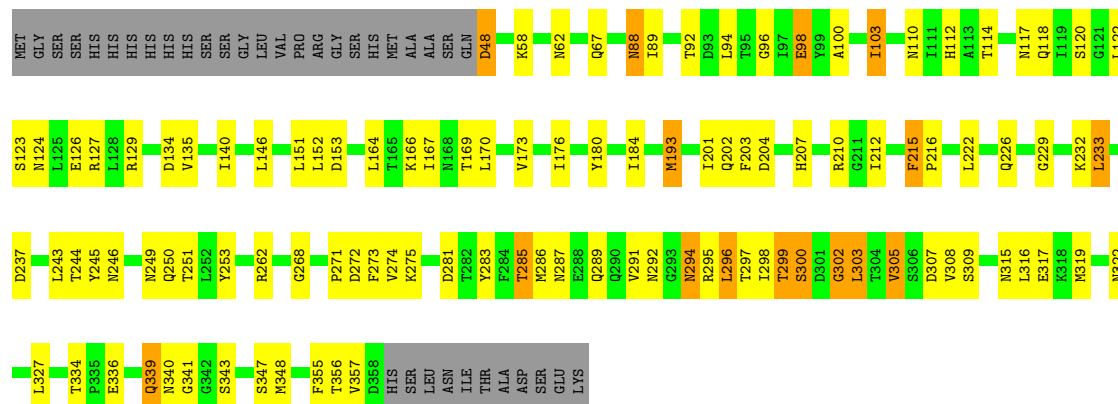
Chain C: 





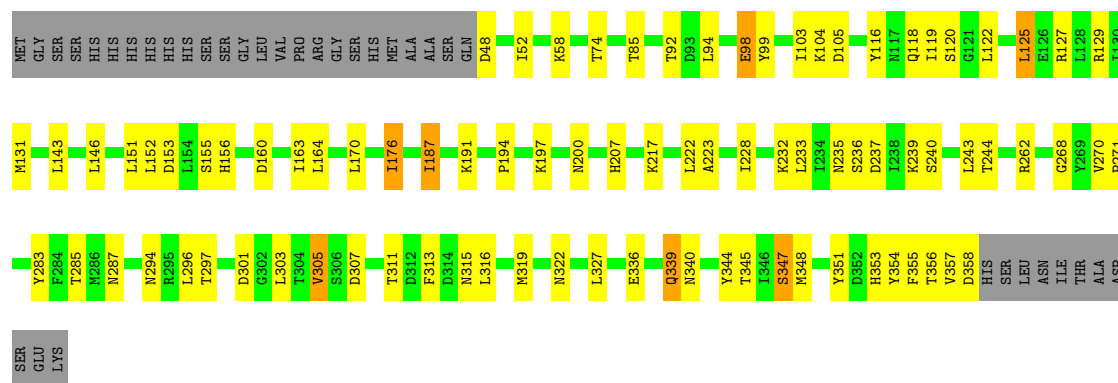
• Molecule 1: Internalin K

Chain D: 58% 27% 5% 10%



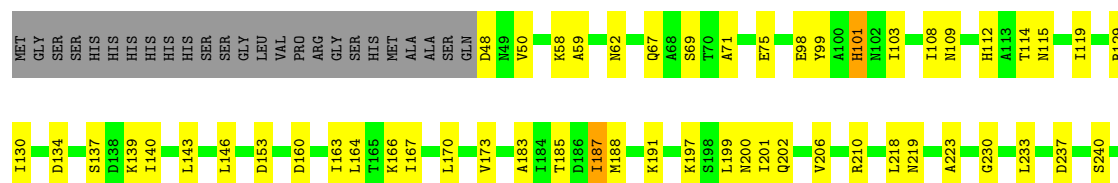
• Molecule 1: Internalin K

Chain E: 65% 23% 10%



• Molecule 1: Internalin K

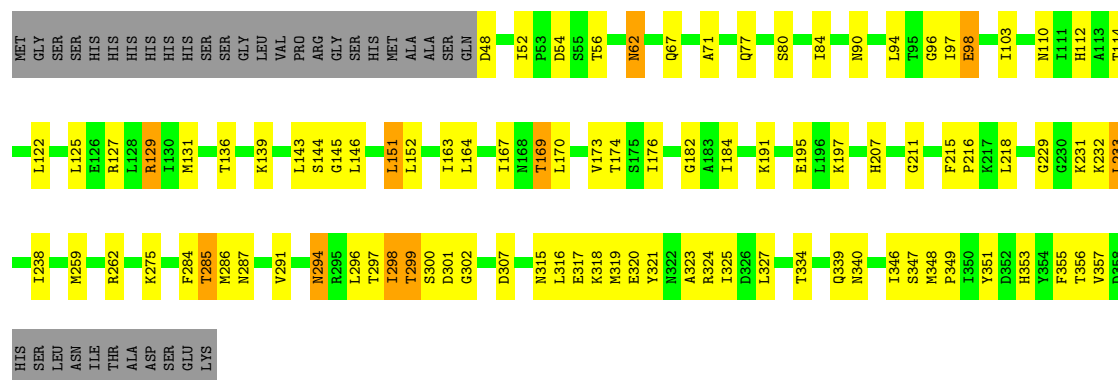
Chain F: 64% 24% 10%





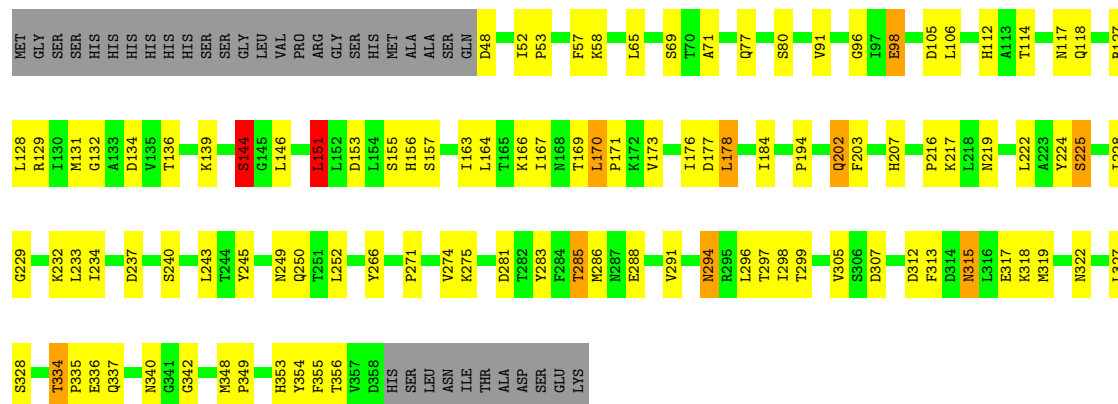
• Molecule 1: Internalin K

Chain G: 62% 24% 10%



• Molecule 1: Internalin K

Chain H: 59% 27% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.10Å 89.50Å 186.30Å 90.00° 103.50° 90.00°	Depositor
Resolution (Å)	47.90 – 2.39 47.90 – 2.40	Depositor EDS
% Data completeness (in resolution range)	80.7 (47.90-2.39) 80.8 (47.90-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.192 , 0.232 0.193 , 0.233	Depositor DCC
R_{free} test set	9215 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	1.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 17.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.35$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	0.299 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.679 for H, K, L 0.321 for -H, -K, H+L	Depositor
Outliers	0 of 92214 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20171	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6461e-03.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NA, NI

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	1.03	1/2478 (0.0%)	1.15	5/3373 (0.1%)
1	B	1.17	9/2478 (0.4%)	1.23	13/3373 (0.4%)
1	C	1.10	3/2478 (0.1%)	1.18	6/3373 (0.2%)
1	D	1.04	4/2478 (0.2%)	1.20	19/3373 (0.6%)
1	E	0.92	0/2478	1.10	8/3373 (0.2%)
1	F	1.10	3/2478 (0.1%)	1.22	15/3373 (0.4%)
1	G	1.02	0/2478	1.20	14/3373 (0.4%)
1	H	1.09	5/2478 (0.2%)	1.21	12/3373 (0.4%)
All	All	1.06	25/19824 (0.1%)	1.19	92/26984 (0.3%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	101	HIS	CG-ND1	-8.21	1.29	1.38
1	B	156	HIS	CG-ND1	-8.11	1.29	1.38
1	F	101	HIS	CD2-NE2	-6.90	1.30	1.37
1	B	156	HIS	CD2-NE2	-6.38	1.30	1.37
1	D	152	LEU	CA-C	6.31	1.60	1.52
1	C	222	LEU	CA-C	-6.11	1.45	1.53
1	B	112	HIS	CG-CD2	5.96	1.42	1.35
1	F	343	SER	CA-C	-5.84	1.45	1.52
1	B	199	LEU	CA-C	5.74	1.59	1.52
1	D	343	SER	CA-C	-5.73	1.45	1.52
1	C	111	ILE	C-O	-5.67	1.18	1.24
1	H	222	LEU	C-O	-5.53	1.17	1.23
1	A	151	LEU	C-O	-5.46	1.17	1.23
1	H	348	MET	C-O	-5.41	1.21	1.23
1	B	354	TYR	N-CA	5.36	1.52	1.45
1	H	207	HIS	CG-CD2	5.21	1.41	1.35
1	B	126	GLU	N-CA	5.20	1.52	1.46
1	H	202	GLN	C-O	-5.18	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	156	HIS	ND1-CE1	5.17	1.37	1.32
1	D	103	ILE	CA-CB	5.17	1.59	1.54
1	D	207	HIS	CG-CD2	5.15	1.41	1.35
1	C	151	LEU	N-CA	5.14	1.52	1.46
1	B	201	ILE	CA-C	5.12	1.59	1.52
1	B	112	HIS	CG-ND1	-5.12	1.32	1.38
1	B	343	SER	C-O	-5.11	1.17	1.23

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	52	ILE	CA-C-N	8.45	128.57	119.28
1	H	52	ILE	C-N-CA	8.45	128.57	119.28
1	C	170	LEU	CA-C-N	7.77	127.83	119.28
1	C	170	LEU	C-N-CA	7.77	127.83	119.28
1	H	144	SER	N-CA-C	7.63	120.56	111.33
1	F	340	ASN	N-CA-C	7.47	121.98	111.30
1	D	58	LYS	N-CA-C	7.39	120.28	111.33
1	B	117	ASN	N-CA-C	7.36	118.94	111.07
1	G	197	LYS	N-CA-C	-7.12	104.46	113.43
1	B	348	MET	CA-C-N	-7.00	113.59	120.52
1	B	348	MET	C-N-CA	-7.00	113.59	120.52
1	D	167	ILE	N-CA-C	-6.88	106.59	113.20
1	G	62	ASN	N-CA-C	-6.88	103.83	111.82
1	D	215	PHE	CA-C-N	6.87	126.33	118.85
1	D	215	PHE	C-N-CA	6.87	126.33	118.85
1	E	170	LEU	CA-C-N	6.73	126.32	119.19
1	E	170	LEU	C-N-CA	6.73	126.32	119.19
1	G	52	ILE	CA-C-N	6.67	126.92	119.32
1	G	52	ILE	C-N-CA	6.67	126.92	119.32
1	G	238	ILE	N-CA-C	6.61	118.31	108.46
1	F	302	GLY	N-CA-C	6.49	119.95	110.80
1	F	140	ILE	CA-C-N	6.37	126.33	120.03
1	F	140	ILE	C-N-CA	6.37	126.33	120.03
1	F	336	GLU	N-CA-C	6.33	118.99	111.33
1	F	130	ILE	N-CA-C	-6.27	98.49	107.77
1	B	193	MET	CA-C-N	-6.26	112.01	119.84
1	B	193	MET	C-N-CA	-6.26	112.01	119.84
1	A	103	ILE	N-CA-CB	6.25	117.99	110.31
1	G	339	GLN	N-CA-C	6.25	119.92	112.93
1	E	85	THR	N-CA-C	6.11	118.89	108.02
1	H	117	ASN	N-CA-C	6.07	118.69	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	187	ILE	CB-CA-C	-6.07	104.15	111.85
1	H	296	LEU	N-CA-C	6.03	119.30	109.59
1	D	302	GLY	N-CA-C	6.00	120.39	110.55
1	B	340	ASN	N-CA-C	5.97	119.84	111.30
1	C	139	LYS	CA-C-N	-5.96	116.88	123.43
1	C	139	LYS	C-N-CA	-5.96	116.88	123.43
1	A	59	ALA	N-CA-C	-5.93	104.89	111.36
1	F	109	ASN	N-CA-C	5.87	118.28	108.02
1	H	348	MET	CA-C-N	-5.86	114.72	120.52
1	H	348	MET	C-N-CA	-5.86	114.72	120.52
1	H	139	LYS	N-CA-C	-5.84	106.15	113.28
1	G	302	GLY	N-CA-C	5.78	119.18	110.75
1	D	341	GLY	N-CA-C	-5.76	106.70	115.00
1	G	348	MET	CA-C-N	-5.71	114.72	120.31
1	G	348	MET	C-N-CA	-5.71	114.72	120.31
1	D	110	ASN	CA-C-N	-5.65	115.13	121.85
1	D	110	ASN	C-N-CA	-5.65	115.13	121.85
1	E	339	GLN	CA-C-N	-5.64	113.71	123.25
1	E	339	GLN	C-N-CA	-5.64	113.71	123.25
1	D	339	GLN	CA-C-N	-5.62	113.63	122.79
1	D	339	GLN	C-N-CA	-5.62	113.63	122.79
1	A	159	HIS	N-CA-C	5.62	118.42	110.50
1	G	211	GLY	N-CA-C	-5.58	107.83	114.48
1	E	103	ILE	N-CA-CB	5.57	116.60	110.53
1	A	348	MET	CA-C-N	-5.56	115.03	120.98
1	A	348	MET	C-N-CA	-5.56	115.03	120.98
1	C	348	MET	CA-C-N	-5.55	114.61	120.66
1	C	348	MET	C-N-CA	-5.55	114.61	120.66
1	D	89	ILE	CA-C-N	-5.52	114.60	122.77
1	D	89	ILE	C-N-CA	-5.52	114.60	122.77
1	H	112	HIS	N-CA-C	5.50	119.17	111.74
1	F	257	SER	N-CA-C	-5.50	106.53	113.18
1	G	339	GLN	CB-CA-C	-5.50	101.30	110.81
1	F	108	ILE	N-CA-C	5.44	115.69	107.75
1	D	120	SER	N-CA-C	5.38	119.34	112.34
1	F	329	TYR	N-CA-C	5.35	117.80	111.33
1	F	330	GLN	CA-C-N	5.30	130.42	121.14
1	F	330	GLN	C-N-CA	5.30	130.42	121.14
1	H	166	LYS	N-CA-C	-5.28	106.84	113.28
1	B	69	SER	N-CA-C	5.25	117.69	111.33
1	F	230	GLY	N-CA-C	5.21	119.44	112.77
1	E	340	ASN	N-CA-C	5.21	118.75	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	296	LEU	N-CA-C	5.21	117.89	109.40
1	F	59	ALA	N-CA-C	-5.17	105.64	111.28
1	B	334	THR	CA-C-N	-5.16	114.92	120.03
1	B	334	THR	C-N-CA	-5.16	114.92	120.03
1	D	135	VAL	N-CA-C	5.16	114.70	107.37
1	D	193	MET	CA-C-N	-5.15	113.40	119.84
1	D	193	MET	C-N-CA	-5.15	113.40	119.84
1	D	296	LEU	N-CA-C	5.14	117.78	109.40
1	D	348	MET	CA-C-N	-5.14	115.30	120.85
1	D	348	MET	C-N-CA	-5.14	115.30	120.85
1	G	184	ILE	CB-CA-C	5.13	117.17	110.96
1	B	129	ARG	CG-CD-NE	-5.10	100.78	112.00
1	B	161	ASP	N-CA-C	5.09	118.26	111.75
1	H	336	GLU	N-CA-C	5.08	117.48	111.33
1	B	215	PHE	CA-C-N	-5.02	113.56	119.84
1	B	215	PHE	C-N-CA	-5.02	113.56	119.84
1	G	151	LEU	N-CA-C	5.02	116.58	108.34
1	H	151	LEU	N-CA-C	5.01	117.07	108.90
1	F	357	VAL	N-CA-C	5.01	115.79	108.53

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	2436	0	2375	53	0
1	B	2436	0	2375	82	0
1	C	2436	0	2375	83	0
1	D	2436	0	2375	91	0
1	E	2436	0	2375	63	0
1	F	2436	0	2375	61	0
1	G	2436	0	2375	67	1
1	H	2436	0	2375	71	0
2	B	1	0	0	0	0
2	F	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	96	0	0	15	0
4	B	93	0	0	29	0
4	C	99	0	0	25	0
4	D	95	0	0	38	0
4	E	59	0	0	17	0
4	F	87	0	0	15	0
4	G	65	0	0	15	0
4	H	85	0	0	19	1
All	All	20171	0	19000	558	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 (558) 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:H:217:LYS:HG3	4:H:579:HOH:O	1.35	1.26
1:C:328:SER:HB3	4:C:486:HOH:O	1.19	1.25
1:B:269:TYR:HA	4:B:574:HOH:O	1.32	1.23
1:B:239:LYS:NZ	4:B:539:HOH:O	1.71	1.22
1:D:298:ILE:HB	4:D:492:HOH:O	1.43	1.15
1:F:345:THR:HB	4:F:576:HOH:O	1.46	1.15
1:A:334:THR:HG23	4:A:422:HOH:O	1.43	1.15
1:B:338:PHE:CD2	4:B:581:HOH:O	2.00	1.14
1:B:285:THR:HG21	4:B:583:HOH:O	1.48	1.11
1:F:328:SER:O	1:F:331:SER:HB3	1.52	1.09
1:C:217:LYS:HE3	4:C:494:HOH:O	1.53	1.09
1:E:160:ASP:O	1:E:163:ILE:HG22	1.56	1.06
1:H:128:LEU:CD1	4:H:584:HOH:O	2.04	1.05
1:F:139:LYS:HD3	4:F:581:HOH:O	1.60	1.02
1:H:106:LEU:HG	4:H:584:HOH:O	1.60	1.00
1:C:166:LYS:HE2	4:C:487:HOH:O	1.61	1.00
1:C:131:MET:HE2	4:C:484:HOH:O	1.61	0.98
1:A:104:LYS:HE3	4:A:479:HOH:O	1.64	0.97
1:C:104:LYS:NZ	4:C:457:HOH:O	1.98	0.96
1:E:345:THR:HG23	4:E:456:HOH:O	1.63	0.96
1:A:261:GLU:O	1:A:351:TYR:HE1	1.49	0.96
1:H:337:GLN:HG2	4:H:572:HOH:O	1.65	0.96
1:D:272:ASP:HA	4:D:451:HOH:O	1.65	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:MET:HG2	4:D:494:HOH:O	1.65	0.94
1:C:237:ASP:OD2	4:C:406:HOH:O	1.87	0.93
1:H:334:THR:HG23	4:H:567:HOH:O	1.70	0.91
1:C:129:ARG:HB3	4:C:484:HOH:O	1.69	0.91
1:B:340:ASN:ND2	4:B:543:HOH:O	2.03	0.91
1:G:207:HIS:HD2	4:G:555:HOH:O	1.51	0.91
1:F:137:SER:HB3	4:F:585:HOH:O	1.71	0.91
1:B:83:TYR:HE1	4:B:586:HOH:O	1.52	0.90
1:D:273:PHE:N	4:D:451:HOH:O	2.06	0.89
1:H:106:LEU:CG	4:H:584:HOH:O	2.19	0.89
1:E:336:GLU:O	4:E:442:HOH:O	1.91	0.88
1:B:73:ILE:HB	4:B:585:HOH:O	1.75	0.87
1:G:54:ASP:HA	4:G:565:HOH:O	1.73	0.86
1:B:93:ASP:OD2	4:B:538:HOH:O	1.94	0.84
1:E:239:LYS:HD3	1:E:358:ASP:HB3	1.59	0.84
1:E:99:TYR:OH	4:E:433:HOH:O	1.95	0.83
1:B:328:SER:HA	1:B:348:MET:HE2	1.60	0.83
1:A:261:GLU:O	1:A:351:TYR:CE1	2.31	0.83
1:D:262:ARG:NH2	4:D:451:HOH:O	2.13	0.82
1:F:139:LYS:HB3	4:F:581:HOH:O	1.80	0.81
1:A:285:THR:CG2	4:A:492:HOH:O	2.28	0.80
1:C:334:THR:HG23	4:C:417:HOH:O	1.83	0.79
1:F:237:ASP:OD1	1:F:356:THR:OG1	2.01	0.79
1:B:48:ASP:OD1	1:B:48:ASP:O	2.01	0.79
1:B:117:ASN:HB3	4:C:480:HOH:O	1.83	0.78
1:C:311:THR:HB	4:C:492:HOH:O	1.85	0.77
1:C:329:TYR:HE1	4:C:477:HOH:O	1.66	0.77
1:F:163:ILE:HG23	1:F:164:LEU:HD12	1.67	0.77
1:C:217:LYS:HG2	4:C:476:HOH:O	1.83	0.77
1:H:216:PRO:HB2	4:H:579:HOH:O	1.85	0.76
1:F:115:ASN:HA	4:F:581:HOH:O	1.85	0.76
1:C:239:LYS:HG3	1:E:301:ASP:OD2	1.85	0.76
1:D:272:ASP:C	4:D:451:HOH:O	2.27	0.76
1:A:48:ASP:OD1	1:A:48:ASP:O	2.04	0.76
1:G:207:HIS:CD2	4:G:555:HOH:O	2.33	0.75
1:B:269:TYR:HD1	4:B:574:HOH:O	1.68	0.75
1:B:109:ASN:OD1	4:B:524:HOH:O	2.05	0.74
1:A:308:VAL:HB	4:A:495:HOH:O	1.86	0.74
1:B:175:SER:HB2	4:B:572:HOH:O	1.88	0.74
1:E:197:LYS:HE2	4:E:434:HOH:O	1.86	0.74
1:D:299:THR:OG1	1:D:300:SER:N	2.15	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:340:ASN:HB2	4:D:474:HOH:O	1.87	0.73
1:E:127:ARG:HG2	1:E:151:LEU:HD22	1.71	0.73
1:C:122:LEU:HB2	1:C:125:LEU:HD22	1.71	0.72
1:A:96:GLY:N	1:A:98:GLU:OE1	2.22	0.72
1:E:313:PHE:HZ	4:E:459:HOH:O	1.71	0.72
1:B:269:TYR:CD1	4:B:574:HOH:O	2.43	0.71
1:C:58:LYS:HE2	1:C:71:ALA:O	1.89	0.71
1:D:124:ASN:HB2	4:D:408:HOH:O	1.91	0.71
1:B:204:ASP:C	1:B:226:GLN:OE1	2.34	0.70
1:B:98:GLU:HG3	1:B:118:GLN:HB3	1.73	0.70
1:D:340:ASN:CB	4:D:474:HOH:O	2.39	0.70
1:D:48:ASP:OD2	4:D:406:HOH:O	2.09	0.70
1:F:202:GLN:NE2	4:F:537:HOH:O	2.19	0.70
1:D:286:MET:HG2	4:D:494:HOH:O	1.90	0.69
1:H:312:ASP:OD2	4:H:532:HOH:O	2.08	0.69
1:F:185:THR:HG22	4:F:563:HOH:O	1.92	0.69
1:B:53:PRO:HD2	4:B:532:HOH:O	1.91	0.69
1:D:275:LYS:HA	4:D:492:HOH:O	1.93	0.69
1:C:90:ASN:HB2	4:C:481:HOH:O	1.93	0.69
1:F:69:SER:O	4:F:520:HOH:O	2.08	0.69
1:C:329:TYR:CE1	4:C:477:HOH:O	2.42	0.69
1:E:339:GLN:O	4:E:401:HOH:O	2.11	0.68
1:D:202:GLN:O	1:D:203:PHE:HB2	1.93	0.68
1:B:220:THR:HG23	1:B:343:SER:HB2	1.74	0.68
1:C:101:HIS:HD2	4:C:485:HOH:O	1.75	0.68
1:D:243:LEU:HD21	1:D:319:MET:HE1	1.76	0.68
1:E:120:SER:O	4:E:404:HOH:O	2.12	0.68
1:G:340:ASN:HB2	4:G:557:HOH:O	1.94	0.68
1:F:334:THR:HG23	4:F:568:HOH:O	1.93	0.67
1:G:340:ASN:CG	4:G:557:HOH:O	2.35	0.67
1:B:204:ASP:O	1:B:226:GLN:OE1	2.13	0.67
1:F:200:ASN:OD1	1:F:202:GLN:HG3	1.94	0.67
1:D:336:GLU:HG2	4:D:486:HOH:O	1.93	0.67
1:E:207:HIS:CD2	4:E:453:HOH:O	2.48	0.66
1:A:239:LYS:HD3	1:A:358:ASP:HB3	1.77	0.66
1:A:285:THR:HG22	4:A:492:HOH:O	1.89	0.66
1:D:262:ARG:NH2	1:D:271:PRO:O	2.28	0.66
1:B:160:ASP:O	1:B:163:ILE:HG22	1.95	0.66
1:G:48:ASP:N	4:G:539:HOH:O	2.27	0.66
1:F:67:GLN:HG2	4:F:566:HOH:O	1.94	0.66
1:G:321:TYR:CE1	1:G:323:ALA:HB2	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:ASP:OD1	1:E:48:ASP:O	2.14	0.65
1:A:294:ASN:OD1	1:A:294:ASN:N	2.29	0.65
1:E:116:TYR:O	1:E:119:ILE:HG12	1.97	0.65
1:C:271:PRO:HG3	1:C:327:LEU:HD13	1.77	0.65
1:H:106:LEU:CB	4:H:584:HOH:O	2.44	0.65
1:H:294:ASN:OD1	1:H:294:ASN:N	2.27	0.65
1:D:176:ILE:HD11	1:D:193:MET:HE1	1.78	0.64
1:F:328:SER:O	1:F:331:SER:CB	2.37	0.64
1:E:232:LYS:HA	4:E:457:HOH:O	1.97	0.64
1:A:160:ASP:O	1:A:163:ILE:HG22	1.98	0.64
1:B:120:SER:OG	1:C:142:ASN:CG	2.41	0.64
1:E:287:ASN:HD21	1:E:316:LEU:HA	1.61	0.64
1:A:306:SER:HB3	4:A:493:HOH:O	1.97	0.64
1:D:272:ASP:CA	4:D:451:HOH:O	2.30	0.63
1:E:347:SER:HB2	4:E:455:HOH:O	1.97	0.63
1:C:247:ALA:HB2	4:C:482:HOH:O	1.99	0.63
1:D:285:THR:HG22	4:D:491:HOH:O	1.98	0.63
1:D:285:THR:CG2	4:D:491:HOH:O	2.46	0.63
1:G:110:ASN:OD1	1:G:112:HIS:CE1	2.51	0.63
1:G:340:ASN:CB	4:G:557:HOH:O	2.46	0.62
1:E:156:HIS:HE1	4:E:451:HOH:O	1.81	0.62
1:C:129:ARG:NH2	1:D:129:ARG:NE	2.48	0.62
1:E:236:SER:HB2	1:E:355:PHE:CD1	2.35	0.62
1:E:262:ARG:NH2	1:E:271:PRO:O	2.33	0.61
1:F:62:ASN:OD1	1:F:67:GLN:NE2	2.32	0.61
1:B:161:ASP:O	1:B:164:LEU:HB2	2.01	0.61
1:B:120:SER:OG	1:C:142:ASN:CB	2.49	0.61
1:B:334:THR:HG21	4:B:502:HOH:O	2.01	0.61
1:C:237:ASP:OD1	1:C:356:THR:OG1	2.14	0.61
1:E:127:ARG:NH1	4:E:418:HOH:O	2.32	0.61
1:D:122:LEU:O	1:D:124:ASN:N	2.34	0.61
1:G:122:LEU:O	1:G:125:LEU:HB2	2.01	0.60
1:H:224:TYR:CD2	1:H:225:SER:HB2	2.36	0.60
1:C:179:SER:O	1:C:180:TYR:HB2	2.00	0.60
1:D:202:GLN:HE21	1:D:222:LEU:HB3	1.65	0.60
1:B:215:PHE:HB3	4:B:570:HOH:O	2.01	0.60
1:B:262:ARG:NH2	1:B:271:PRO:O	2.33	0.60
1:H:202:GLN:O	1:H:203:PHE:HB2	2.02	0.60
1:C:246:ASN:OD1	1:C:249:ASN:HB2	2.02	0.60
1:A:261:GLU:C	1:A:351:TYR:HE1	2.09	0.59
1:C:286:MET:HE2	1:C:291:VAL:HG22	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:183:ALA:HA	4:F:507:HOH:O	2.02	0.59
1:F:319:MET:HB2	1:F:355:PHE:HB2	1.83	0.59
1:A:183:ALA:HB2	1:A:231:LYS:HB2	1.83	0.59
1:G:325:ILE:HB	1:G:349:PRO:HG2	1.85	0.59
1:H:252:LEU:HB3	1:H:305:VAL:HB	1.84	0.59
1:A:328:SER:HB2	4:A:416:HOH:O	2.02	0.59
1:F:271:PRO:HG3	1:F:327:LEU:HD13	1.84	0.59
1:B:337:GLN:HG3	4:B:581:HOH:O	2.04	0.58
1:C:201:ILE:HD12	1:C:206:VAL:HG21	1.84	0.58
1:G:96:GLY:N	1:G:98:GLU:OE1	2.31	0.58
1:G:275:LYS:HD3	1:G:298:ILE:HG22	1.84	0.58
1:D:283:TYR:HB3	1:D:322:ASN:HB3	1.85	0.58
1:A:129:ARG:NH2	1:B:129:ARG:NH2	2.51	0.58
1:A:271:PRO:HG3	1:A:327:LEU:HD13	1.83	0.58
1:D:340:ASN:CG	4:D:474:HOH:O	2.46	0.58
1:E:223:ALA:HB3	1:E:345:THR:O	2.03	0.58
1:A:104:LYS:HG3	4:A:490:HOH:O	2.03	0.58
1:F:129:ARG:HG2	1:F:153:ASP:HB3	1.85	0.58
1:C:153:ASP:OD1	1:C:155:SER:OG	2.15	0.57
1:F:188:MET:HE3	1:F:210:ARG:O	2.04	0.57
1:C:129:ARG:NH2	1:D:129:ARG:HE	2.02	0.57
1:G:129:ARG:HD3	1:G:131:MET:HE2	1.87	0.57
1:G:163:ILE:HG23	1:G:164:LEU:HD12	1.86	0.57
1:A:237:ASP:OD1	1:A:356:THR:OG1	2.20	0.57
1:A:261:GLU:C	1:A:351:TYR:CE1	2.83	0.57
1:F:48:ASP:N	4:F:549:HOH:O	2.38	0.57
1:G:315:ASN:HB3	4:G:556:HOH:O	2.05	0.56
1:D:180:TYR:CZ	1:D:203:PHE:CE2	2.93	0.56
1:E:98:GLU:CG	1:E:118:GLN:HB3	2.35	0.56
1:C:250:GLN:NE2	1:C:309:SER:HA	2.21	0.56
1:C:319:MET:HB2	1:C:355:PHE:HB2	1.86	0.56
1:C:197:LYS:O	1:C:218:LEU:HD12	2.06	0.56
1:C:287:ASN:HA	1:C:318:LYS:HB2	1.88	0.56
1:D:134:ASP:CA	4:D:479:HOH:O	2.54	0.56
1:C:48:ASP:HB3	4:C:474:HOH:O	2.05	0.56
1:D:122:LEU:C	1:D:124:ASN:H	2.12	0.56
1:E:104:LYS:NZ	4:E:426:HOH:O	2.37	0.55
1:G:287:ASN:ND2	1:G:315:ASN:O	2.39	0.55
1:G:317:GLU:O	1:G:356:THR:HA	2.06	0.55
1:C:287:ASN:HD21	1:C:316:LEU:HD12	1.71	0.55
1:D:134:ASP:HA	4:D:479:HOH:O	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:277:THR:HG22	1:F:298:ILE:HD12	1.87	0.55
1:G:77:GLN:HG2	4:G:527:HOH:O	2.07	0.55
1:B:269:TYR:CB	4:B:574:HOH:O	2.51	0.55
1:H:106:LEU:HB3	4:H:584:HOH:O	2.04	0.55
1:G:77:GLN:O	1:G:80:SER:OG	2.16	0.55
1:H:134:ASP:CG	4:H:520:HOH:O	2.49	0.55
1:B:335:PRO:HB2	4:B:581:HOH:O	2.07	0.55
1:B:328:SER:HA	1:B:348:MET:CE	2.33	0.55
1:H:271:PRO:HG3	1:H:327:LEU:HD13	1.89	0.55
1:H:286:MET:HE2	1:H:291:VAL:CG2	2.37	0.55
1:H:48:ASP:N	4:H:507:HOH:O	2.39	0.54
1:D:122:LEU:C	1:D:124:ASN:N	2.63	0.54
1:G:285:THR:HG21	4:G:528:HOH:O	2.06	0.54
1:D:229:GLY:HA2	1:D:232:LYS:O	2.07	0.54
1:B:50:VAL:HB	4:B:585:HOH:O	2.06	0.54
1:B:338:PHE:CE2	4:B:581:HOH:O	2.45	0.54
1:B:338:PHE:HD2	4:B:581:HOH:O	1.57	0.54
1:F:295:ARG:NH2	1:F:308:VAL:HG23	2.23	0.54
1:E:283:TYR:HB3	1:E:322:ASN:HB3	1.90	0.54
1:F:187:ILE:HD13	1:F:199:LEU:HD23	1.90	0.54
1:G:127:ARG:HG2	1:G:151:LEU:HB3	1.88	0.54
1:H:128:LEU:HD12	4:H:584:HOH:O	1.85	0.54
1:C:48:ASP:N	4:C:443:HOH:O	2.40	0.54
1:C:201:ILE:CD1	1:C:206:VAL:HG21	2.38	0.54
1:E:98:GLU:HG3	1:E:118:GLN:HB3	1.88	0.54
1:G:62:ASN:OD1	1:G:67:GLN:NE2	2.40	0.54
1:B:146:LEU:HB3	1:B:149:LEU:HB2	1.90	0.54
1:D:98:GLU:HG3	1:D:118:GLN:HB3	1.89	0.54
1:D:134:ASP:CB	4:D:479:HOH:O	2.56	0.54
1:B:340:ASN:ND2	4:B:504:HOH:O	2.39	0.53
1:E:127:ARG:CZ	4:E:436:HOH:O	2.55	0.53
1:C:246:ASN:ND2	4:C:452:HOH:O	2.40	0.53
1:F:112:HIS:HA	1:F:134:ASP:OD2	2.08	0.53
1:F:210:ARG:HA	4:F:534:HOH:O	2.08	0.53
1:B:153:ASP:HB2	4:B:591:HOH:O	2.08	0.53
1:E:351:TYR:CE2	1:E:353:HIS:HB2	2.44	0.53
1:F:163:ILE:O	1:F:166:LYS:HB2	2.09	0.53
1:F:271:PRO:HB3	1:F:327:LEU:HD13	1.89	0.53
1:F:308:VAL:HG13	1:F:312:ASP:HB2	1.91	0.53
1:D:134:ASP:HB3	4:D:479:HOH:O	2.09	0.53
1:H:58:LYS:HE2	1:H:71:ALA:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:THR:O	1:A:139:LYS:HE2	2.08	0.53
1:A:152:LEU:O	1:A:176:ILE:HA	2.09	0.53
1:A:75:GLU:O	4:A:445:HOH:O	2.19	0.52
1:D:246:ASN:OD1	1:D:249:ASN:N	2.28	0.52
1:E:311:THR:HG22	1:E:315:ASN:ND2	2.24	0.52
1:H:98:GLU:HG3	1:H:118:GLN:HB3	1.92	0.52
1:F:340:ASN:CG	1:F:340:ASN:O	2.46	0.52
1:F:163:ILE:CG2	1:F:164:LEU:HD12	2.39	0.52
1:C:141:PRO:HA	4:C:487:HOH:O	2.09	0.52
1:G:229:GLY:O	1:G:233:LEU:HD12	2.10	0.52
1:A:351:TYR:CE2	1:A:353:HIS:HB2	2.45	0.52
1:E:207:HIS:HB3	1:E:228:ILE:HG12	1.91	0.52
1:F:170:LEU:HB3	1:F:173:VAL:CG2	2.40	0.52
1:D:96:GLY:N	1:D:98:GLU:OE1	2.38	0.51
1:D:245:TYR:HD2	4:D:464:HOH:O	1.92	0.51
1:E:319:MET:HB2	1:E:355:PHE:HB2	1.92	0.51
1:F:294:ASN:OD1	1:F:294:ASN:N	2.43	0.51
1:C:275:LYS:HD3	1:C:298:ILE:HG22	1.92	0.51
1:C:287:ASN:ND2	1:C:316:LEU:HD12	2.25	0.51
1:C:287:ASN:HD21	1:C:316:LEU:HA	1.75	0.51
1:F:201:ILE:HD13	1:F:206:VAL:HG21	1.92	0.51
1:G:294:ASN:N	1:G:294:ASN:OD1	2.44	0.51
1:D:233:LEU:HB2	4:D:445:HOH:O	2.10	0.51
1:E:131:MET:HG2	1:E:155:SER:O	2.11	0.51
1:F:160:ASP:O	1:F:163:ILE:HG22	2.10	0.51
1:F:170:LEU:HB3	1:F:173:VAL:HG23	1.91	0.51
1:G:259:MET:HE3	1:G:321:TYR:CD2	2.46	0.51
1:H:48:ASP:OD1	1:H:48:ASP:O	2.29	0.51
1:H:194:PRO:HD2	4:H:514:HOH:O	2.11	0.51
1:H:219:ASN:OD1	1:H:219:ASN:C	2.54	0.51
1:B:177:ASP:OD2	4:B:509:HOH:O	2.19	0.51
1:D:48:ASP:CB	4:D:406:HOH:O	2.58	0.51
1:D:216:PRO:HD3	4:D:418:HOH:O	2.11	0.51
1:E:237:ASP:OD1	1:E:356:THR:OG1	2.27	0.51
1:H:170:LEU:HG	1:H:173:VAL:HG21	1.92	0.51
1:C:247:ALA:CB	4:C:482:HOH:O	2.58	0.51
1:C:176:ILE:HD11	1:C:193:MET:HE1	1.94	0.50
1:G:351:TYR:CE2	1:G:353:HIS:HB2	2.46	0.50
1:E:311:THR:HG22	1:E:315:ASN:HD21	1.75	0.50
1:D:210:ARG:HD3	4:D:481:HOH:O	2.11	0.50
1:F:129:ARG:HG3	1:F:129:ARG:HH11	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:TYR:O	1:B:119:ILE:HG12	2.11	0.50
1:D:286:MET:HG3	1:D:291:VAL:CG2	2.41	0.50
1:D:317:GLU:O	1:D:356:THR:HA	2.11	0.50
1:E:194:PRO:O	1:E:217:LYS:HD2	2.11	0.50
1:A:189:PRO:HD3	4:A:474:HOH:O	2.11	0.50
1:B:143:LEU:HB3	1:B:146:LEU:HD22	1.92	0.50
1:D:124:ASN:OD1	4:D:484:HOH:O	2.19	0.50
1:E:92:THR:O	1:E:92:THR:HG22	2.11	0.50
1:F:187:ILE:HG21	1:F:201:ILE:HG21	1.94	0.50
1:G:262:ARG:HB2	4:G:559:HOH:O	2.12	0.50
1:B:244:THR:HB	1:B:253:TYR:HB3	1.92	0.50
1:B:294:ASN:N	1:B:294:ASN:OD1	2.45	0.50
1:E:235:ASN:HA	1:E:354:TYR:O	2.12	0.50
1:D:117:ASN:HB3	4:D:462:HOH:O	2.12	0.49
1:G:298:ILE:HB	4:G:521:HOH:O	2.12	0.49
1:H:170:LEU:HB3	1:H:173:VAL:HG23	1.94	0.49
1:C:186:ASP:C	1:C:186:ASP:OD1	2.54	0.49
1:D:233:LEU:O	4:D:469:HOH:O	2.20	0.49
1:D:268:GLY:HA2	4:D:481:HOH:O	2.11	0.49
1:E:163:ILE:HG23	1:E:164:LEU:HD12	1.94	0.49
1:D:153:ASP:OD1	1:D:153:ASP:C	2.54	0.49
1:H:128:LEU:HD13	4:H:584:HOH:O	1.91	0.49
1:B:79:ASP:O	1:B:102:ASN:ND2	2.43	0.49
1:B:175:SER:HB3	1:B:198:SER:HB3	1.94	0.49
1:D:210:ARG:CG	4:D:481:HOH:O	2.60	0.49
1:D:286:MET:HA	4:D:494:HOH:O	2.12	0.49
1:B:112:HIS:HA	1:B:134:ASP:OD2	2.13	0.49
1:H:127:ARG:HG2	1:H:151:LEU:CB	2.42	0.49
1:A:88:ASN:OD1	1:A:110:ASN:ND2	2.41	0.49
1:D:126:GLU:HB3	4:D:415:HOH:O	2.12	0.49
1:A:295:ARG:O	1:A:305:VAL:HA	2.12	0.49
1:C:87:ALA:HB3	1:D:88:ASN:HB2	1.95	0.49
1:G:163:ILE:CG2	1:G:164:LEU:HD12	2.42	0.49
1:H:127:ARG:HG2	1:H:151:LEU:HB3	1.94	0.49
1:A:235:ASN:HA	1:A:354:TYR:O	2.12	0.48
1:B:340:ASN:CG	4:B:543:HOH:O	2.45	0.48
1:F:143:LEU:HB3	1:F:146:LEU:HD22	1.94	0.48
1:B:148:ASN:ND2	4:B:565:HOH:O	2.46	0.48
1:G:327:LEU:HD23	1:G:346:ILE:O	2.13	0.48
1:A:182:GLY:HA3	1:A:231:LYS:HG3	1.95	0.48
1:E:143:LEU:HB3	1:E:146:LEU:HD22	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:202:GLN:HB3	1:H:203:PHE:CD2	2.48	0.48
1:A:128:LEU:HD23	1:A:152:LEU:HD13	1.94	0.48
1:C:194:PRO:HB2	1:C:195:GLU:HG2	1.93	0.48
1:D:296:LEU:HD22	1:D:303:LEU:HD11	1.95	0.48
1:G:182:GLY:HA3	1:G:231:LYS:HG3	1.96	0.48
1:C:207:HIS:O	1:C:263:THR:HA	2.13	0.48
1:H:129:ARG:HG2	1:H:153:ASP:HB3	1.94	0.48
1:D:170:LEU:HB3	1:D:173:VAL:HG23	1.95	0.48
1:E:271:PRO:HG3	1:E:327:LEU:HD13	1.95	0.48
1:B:98:GLU:HG2	1:C:144:SER:HB2	1.94	0.48
1:H:249:ASN:O	1:H:250:GLN:HB2	2.13	0.48
1:B:178:LEU:HD13	1:B:184:ILE:HD13	1.95	0.48
1:D:202:GLN:NE2	1:D:222:LEU:HB3	2.29	0.48
1:F:286:MET:O	1:F:289:GLN:HB2	2.14	0.48
1:G:151:LEU:C	1:G:151:LEU:HD23	2.39	0.48
1:D:100:ALA:C	4:D:480:HOH:O	2.56	0.48
1:A:285:THR:HB	1:A:320:GLU:HB3	1.96	0.47
1:B:204:ASP:O	1:B:226:GLN:HA	2.14	0.47
1:C:129:ARG:NH2	1:D:129:ARG:NH2	2.62	0.47
1:D:62:ASN:OD1	1:D:67:GLN:NE2	2.47	0.47
1:D:339:GLN:O	4:D:403:HOH:O	2.20	0.47
1:H:144:SER:O	4:H:558:HOH:O	2.20	0.47
1:H:176:ILE:HG22	1:H:177:ASP:N	2.29	0.47
1:H:245:TYR:HB2	1:H:313:PHE:CE1	2.49	0.47
1:A:188:MET:O	1:A:190:LEU:N	2.47	0.47
1:C:219:ASN:OD1	1:C:219:ASN:C	2.57	0.47
1:E:296:LEU:HD23	1:E:305:VAL:HG13	1.97	0.47
1:F:58:LYS:HE2	1:F:71:ALA:O	2.14	0.47
1:G:353:HIS:CD2	1:G:355:PHE:CZ	3.03	0.47
1:A:143:LEU:HB3	1:A:146:LEU:HD22	1.97	0.47
1:A:284:PHE:O	1:A:290:GLN:HA	2.15	0.47
1:H:96:GLY:N	1:H:98:GLU:OE1	2.43	0.47
1:D:48:ASP:O	1:D:48:ASP:OD1	2.33	0.47
1:B:105:ASP:OD1	1:B:127:ARG:CZ	2.63	0.47
1:B:337:GLN:CG	4:B:581:HOH:O	2.63	0.47
1:D:103:ILE:O	1:D:124:ASN:O	2.32	0.47
1:D:244:THR:HB	1:D:253:TYR:HB3	1.97	0.47
1:F:143:LEU:HD12	1:F:167:ILE:HG22	1.95	0.47
1:F:197:LYS:C	1:F:218:LEU:HD12	2.40	0.47
1:H:286:MET:HE2	1:H:291:VAL:HG21	1.95	0.47
1:B:129:ARG:HG2	1:B:153:ASP:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:236:SER:HB2	1:E:355:PHE:HD1	1.77	0.47
1:H:203:PHE:HA	1:H:225:SER:O	2.14	0.47
1:H:237:ASP:OD1	1:H:356:THR:OG1	2.26	0.47
1:G:94:LEU:O	1:G:97:ILE:HG22	2.13	0.47
1:H:315:ASN:OD1	1:H:315:ASN:N	2.47	0.47
1:C:129:ARG:NH2	1:D:129:ARG:CZ	2.77	0.46
1:D:204:ASP:C	1:D:226:GLN:OE1	2.58	0.46
1:E:105:ASP:OD1	1:E:127:ARG:HB2	2.15	0.46
1:E:125:LEU:HD12	1:E:125:LEU:HA	1.67	0.46
1:E:344:TYR:N	4:E:456:HOH:O	2.47	0.46
1:C:322:ASN:HA	1:C:351:TYR:O	2.15	0.46
1:B:89:ILE:HG13	1:B:91:VAL:HG23	1.97	0.46
1:C:132:GLY:O	1:C:157:SER:HA	2.15	0.46
1:C:134:ASP:OD1	1:C:134:ASP:N	2.49	0.46
1:D:253:TYR:CE1	1:D:302:GLY:HA3	2.51	0.46
1:G:174:THR:HA	1:G:195:GLU:O	2.16	0.46
1:G:84:ILE:HD12	1:G:103:ILE:HD13	1.97	0.46
1:B:118:GLN:HG3	1:C:169:THR:HG21	1.98	0.46
1:B:202:GLN:O	1:B:203:PHE:HB2	2.15	0.46
1:F:103:ILE:HG23	4:F:506:HOH:O	2.16	0.46
1:H:266:TYR:CD2	1:H:335:PRO:HG3	2.50	0.46
1:A:112:HIS:HA	1:A:134:ASP:OD2	2.16	0.46
1:C:251:THR:HA	1:C:305:VAL:O	2.15	0.46
1:D:274:VAL:HG21	1:D:281:ASP:HB3	1.97	0.46
1:E:152:LEU:HD23	1:E:176:ILE:HG12	1.97	0.46
1:C:137:SER:O	1:C:166:LYS:NZ	2.49	0.46
1:H:65:LEU:HD11	1:H:77:GLN:HB3	1.98	0.46
1:B:311:THR:HG23	4:B:540:HOH:O	2.16	0.45
1:D:92:THR:HG22	1:D:92:THR:O	2.16	0.45
1:E:52:ILE:HB	1:E:58:LYS:HD2	1.98	0.45
1:G:110:ASN:OD1	1:G:112:HIS:HE1	1.98	0.45
1:B:126:GLU:O	1:B:149:LEU:HD12	2.17	0.45
1:B:164:LEU:HD23	1:B:190:LEU:HG	1.97	0.45
1:C:266:TYR:CD2	1:C:335:PRO:HG3	2.51	0.45
1:D:127:ARG:HG2	1:D:151:LEU:HB3	1.99	0.45
1:H:129:ARG:HG3	1:H:129:ARG:HH11	1.80	0.45
1:C:294:ASN:OD1	1:C:294:ASN:N	2.49	0.45
1:D:166:LYS:HG2	4:D:463:HOH:O	2.16	0.45
1:B:268:GLY:O	1:B:270:VAL:HG23	2.16	0.45
1:B:353:HIS:CD2	1:B:355:PHE:CZ	3.05	0.45
1:D:251:THR:HA	1:D:305:VAL:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:295:ARG:HH22	1:F:308:VAL:CG2	2.29	0.45
1:G:136:THR:O	1:G:139:LYS:HB2	2.17	0.45
1:G:229:GLY:HA2	1:G:232:LYS:O	2.17	0.45
1:B:182:GLY:HA3	1:B:231:LYS:HG2	1.99	0.45
1:H:57:PHE:HB2	1:H:91:VAL:HG13	1.98	0.45
1:C:178:LEU:HD23	1:C:178:LEU:HA	1.82	0.45
1:F:219:ASN:O	1:F:342:GLY:HA3	2.17	0.45
1:G:125:LEU:HD12	1:G:125:LEU:HA	1.77	0.45
1:D:275:LYS:CA	4:D:492:HOH:O	2.60	0.44
1:B:187:ILE:HD12	1:B:199:LEU:CD2	2.47	0.44
1:D:204:ASP:O	1:D:226:GLN:HA	2.17	0.44
1:E:94:LEU:HD11	1:E:116:TYR:CD1	2.52	0.44
1:F:48:ASP:OD1	1:F:48:ASP:O	2.34	0.44
1:H:234:ILE:HG21	1:H:353:HIS:CE1	2.52	0.44
1:B:127:ARG:HG2	1:B:151:LEU:HB3	1.99	0.44
1:C:108:ILE:HG23	1:C:111:ILE:HG21	1.99	0.44
1:F:50:VAL:HG22	1:F:75:GLU:HG2	1.98	0.44
1:G:284:PHE:HE2	1:G:286:MET:HG3	1.83	0.44
1:H:288:GLU:OE2	1:H:318:LYS:HE3	2.17	0.44
1:A:96:GLY:CA	1:A:98:GLU:OE1	2.65	0.44
1:B:292:ASN:OD1	1:B:293:GLY:N	2.50	0.44
1:C:149:LEU:HD12	1:C:149:LEU:HA	1.77	0.44
1:F:223:ALA:HB3	1:F:346:ILE:HG22	2.00	0.44
1:G:191:LYS:HA	1:G:215:PHE:CE1	2.52	0.44
1:H:170:LEU:HB3	1:H:173:VAL:CG2	2.48	0.44
1:H:283:TYR:HB3	1:H:322:ASN:HB3	1.99	0.44
1:B:120:SER:HB3	1:B:141:PRO:HG2	1.99	0.44
1:D:287:ASN:HD21	1:D:316:LEU:HA	1.82	0.44
1:E:129:ARG:HA	1:E:153:ASP:HB3	2.00	0.44
1:G:173:VAL:HG11	1:G:176:ILE:HG13	1.99	0.44
1:G:319:MET:HB2	1:G:355:PHE:HB2	1.98	0.44
1:G:299:THR:C	1:G:301:ASP:N	2.75	0.44
1:H:105:ASP:OD1	1:H:127:ARG:CZ	2.65	0.44
1:H:229:GLY:HA2	1:H:232:LYS:O	2.17	0.44
1:D:245:TYR:OH	1:D:308:VAL:O	2.24	0.44
1:D:319:MET:HB2	1:D:355:PHE:HB2	1.99	0.44
1:C:170:LEU:HD13	1:C:173:VAL:HG21	2.00	0.44
1:C:224:TYR:CZ	4:C:414:HOH:O	2.68	0.44
1:G:318:LYS:HA	1:G:355:PHE:O	2.18	0.44
1:B:48:ASP:O	1:B:48:ASP:CG	2.60	0.43
1:H:285:THR:O	1:H:319:MET:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:ARG:O	1:A:263:THR:C	2.61	0.43
1:C:285:THR:HB	1:C:320:GLU:HB3	1.99	0.43
1:D:271:PRO:HB3	1:D:327:LEU:HD13	2.00	0.43
1:B:271:PRO:HB3	1:B:327:LEU:HD13	1.99	0.43
1:H:132:GLY:O	1:H:157:SER:HA	2.18	0.43
1:A:83:TYR:HE1	4:A:457:HOH:O	2.02	0.43
1:C:117:ASN:HA	1:C:141:PRO:HG3	2.00	0.43
1:D:294:ASN:OD1	1:D:294:ASN:N	2.51	0.43
1:G:143:LEU:O	1:G:145:GLY:N	2.52	0.43
1:H:170:LEU:HA	1:H:171:PRO:HD3	1.75	0.43
1:B:146:LEU:HD23	1:B:149:LEU:HD22	2.01	0.43
1:C:328:SER:CB	4:C:486:HOH:O	2.07	0.43
1:E:191:LYS:HE2	1:E:191:LYS:HB3	1.79	0.43
1:F:274:VAL:HG22	1:F:324:ARG:O	2.19	0.43
1:F:286:MET:HG3	1:F:291:VAL:CG2	2.48	0.43
1:G:62:ASN:OD1	1:G:71:ALA:O	2.36	0.43
1:E:122:LEU:O	1:E:125:LEU:HB2	2.19	0.43
1:H:129:ARG:HG3	1:H:129:ARG:NH1	2.33	0.43
1:A:286:MET:O	1:A:287:ASN:C	2.61	0.43
1:D:292:ASN:HB3	1:D:295:ARG:HG3	2.00	0.43
1:G:170:LEU:HD13	1:G:173:VAL:HG21	2.00	0.43
1:H:353:HIS:CD2	1:H:355:PHE:CZ	3.06	0.43
1:C:69:SER:O	4:C:447:HOH:O	2.20	0.43
1:D:250:GLN:NE2	1:D:309:SER:HA	2.33	0.43
1:E:200:ASN:HB2	1:E:222:LEU:HD12	2.01	0.43
1:G:143:LEU:O	1:G:144:SER:C	2.62	0.43
1:H:131:MET:HB3	1:H:155:SER:O	2.18	0.43
1:C:252:LEU:HB3	1:C:305:VAL:HB	1.99	0.43
1:H:228:ILE:HD11	4:H:581:HOH:O	2.19	0.43
1:B:287:ASN:HA	1:B:318:LYS:HB2	2.00	0.42
1:G:287:ASN:HD21	1:G:316:LEU:HA	1.82	0.42
1:H:77:GLN:O	1:H:80:SER:OG	2.29	0.42
1:H:286:MET:HE2	1:H:291:VAL:HG22	2.00	0.42
1:H:317:GLU:O	1:H:356:THR:HA	2.19	0.42
1:A:319:MET:HB2	1:A:355:PHE:HB2	2.01	0.42
1:C:129:ARG:HH22	1:D:129:ARG:HH21	1.68	0.42
1:D:237:ASP:OD1	1:D:356:THR:OG1	2.33	0.42
1:H:202:GLN:O	1:H:203:PHE:CB	2.67	0.42
1:G:316:LEU:HD23	1:G:357:VAL:HG21	2.01	0.42
1:A:207:HIS:HD2	4:A:448:HOH:O	2.01	0.42
1:D:129:ARG:HG3	1:D:129:ARG:HH11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:98:GLU:HG2	1:E:118:GLN:HB3	2.00	0.42
1:F:285:THR:HA	1:F:289:GLN:O	2.20	0.42
1:F:334:THR:CG2	4:F:568:HOH:O	2.60	0.42
1:D:286:MET:O	1:D:289:GLN:HB2	2.18	0.42
1:G:299:THR:O	1:G:300:SER:C	2.61	0.42
1:H:163:ILE:HG23	1:H:164:LEU:HD12	2.00	0.42
1:A:334:THR:CG2	4:A:422:HOH:O	2.28	0.42
1:B:170:LEU:HA	1:B:171:PRO:HD2	1.95	0.42
1:C:229:GLY:HA2	1:C:232:LYS:O	2.20	0.42
1:E:294:ASN:N	1:E:294:ASN:OD1	2.52	0.42
1:E:348:MET:HB2	4:E:455:HOH:O	2.19	0.42
1:F:129:ARG:HG3	1:F:129:ARG:NH1	2.35	0.42
1:G:144:SER:HB3	1:G:169:THR:OG1	2.19	0.42
1:B:327:LEU:HA	1:B:327:LEU:HD12	1.62	0.42
1:F:295:ARG:NH2	1:F:308:VAL:CG2	2.83	0.42
1:G:48:ASP:OD1	1:G:48:ASP:O	2.37	0.42
1:G:152:LEU:HA	1:G:152:LEU:HD12	1.81	0.42
1:H:275:LYS:HD3	1:H:298:ILE:HG22	2.00	0.42
1:A:129:ARG:HB3	1:A:131:MET:HE2	2.00	0.42
1:B:149:LEU:HD12	1:B:149:LEU:HA	1.81	0.42
1:A:221:LEU:HD12	1:A:221:LEU:HA	1.90	0.42
1:E:336:GLU:O	1:E:339:GLN:HB2	2.20	0.42
1:G:285:THR:HB	1:G:320:GLU:HB3	2.01	0.42
1:D:94:LEU:HD23	1:D:94:LEU:HA	1.79	0.42
1:D:210:ARG:CD	4:D:481:HOH:O	2.67	0.42
1:E:243:LEU:HD21	1:E:319:MET:HE1	2.02	0.42
1:G:315:ASN:CB	4:G:556:HOH:O	2.67	0.42
1:H:53:PRO:HG3	4:H:546:HOH:O	2.20	0.42
1:E:236:SER:HB2	1:E:355:PHE:CE1	2.54	0.41
1:F:99:TYR:HA	1:F:101:HIS:CE1	2.55	0.41
1:H:178:LEU:HD12	1:H:178:LEU:HA	1.97	0.41
1:H:327:LEU:HB3	1:H:349:PRO:HD2	2.02	0.41
1:A:88:ASN:ND2	4:A:450:HOH:O	2.51	0.41
1:G:324:ARG:NH1	4:G:504:HOH:O	2.21	0.41
1:A:340:ASN:ND2	4:A:414:HOH:O	2.40	0.41
1:E:120:SER:HA	1:E:146:LEU:HD11	2.02	0.41
1:E:268:GLY:O	1:E:270:VAL:HG23	2.20	0.41
1:A:48:ASP:O	1:A:48:ASP:CG	2.64	0.41
1:C:129:ARG:HH21	1:D:129:ARG:CZ	2.33	0.41
1:C:170:LEU:HB3	1:C:173:VAL:HG23	2.01	0.41
1:G:191:LYS:HA	1:G:215:PHE:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:THR:CG2	4:E:435:HOH:O	2.68	0.41
1:G:94:LEU:HD23	1:G:94:LEU:HA	1.81	0.41
1:H:274:VAL:HG21	1:H:281:ASP:HB3	2.03	0.41
1:A:48:ASP:OD1	1:A:48:ASP:C	2.63	0.41
1:C:68:ALA:O	1:C:69:SER:C	2.60	0.41
1:F:271:PRO:CB	1:F:327:LEU:HD13	2.50	0.41
1:G:286:MET:HE2	1:G:291:VAL:CG2	2.50	0.41
1:B:175:SER:CB	4:B:572:HOH:O	2.59	0.41
1:B:197:LYS:HA	1:B:218:LEU:HA	2.02	0.41
1:C:122:LEU:CB	1:C:125:LEU:HD22	2.47	0.41
1:F:219:ASN:OD1	1:F:219:ASN:C	2.63	0.41
1:B:117:ASN:HA	1:B:141:PRO:HG3	2.02	0.41
1:C:327:LEU:HD23	1:C:346:ILE:O	2.21	0.41
1:E:164:LEU:HD21	1:E:187:ILE:HA	2.03	0.41
1:E:239:LYS:HD3	1:E:239:LYS:HA	1.88	0.41
1:F:187:ILE:HD13	1:F:199:LEU:CD2	2.51	0.41
1:G:233:LEU:HD12	1:G:233:LEU:HA	1.84	0.41
1:H:340:ASN:CG	1:H:340:ASN:O	2.61	0.41
1:B:340:ASN:CG	1:B:340:ASN:O	2.61	0.41
1:D:170:LEU:HD23	1:D:170:LEU:HA	1.88	0.41
1:G:215:PHE:HA	1:G:216:PRO:HD2	1.94	0.41
1:H:354:TYR:HE1	4:H:583:HOH:O	2.03	0.41
1:C:201:ILE:HD12	1:C:206:VAL:CG2	2.51	0.40
1:D:212:ILE:O	1:D:215:PHE:HB2	2.20	0.40
1:G:163:ILE:HD12	1:G:163:ILE:HA	1.93	0.40
1:G:218:LEU:N	4:G:563:HOH:O	2.49	0.40
1:H:176:ILE:CG2	1:H:177:ASP:N	2.84	0.40
1:A:296:LEU:HD22	1:A:303:LEU:HD11	2.03	0.40
1:A:325:ILE:HD11	1:A:351:TYR:HB2	2.02	0.40
1:B:233:LEU:HD12	1:B:233:LEU:HA	1.84	0.40
1:B:332:TYR:CE1	1:B:346:ILE:HG12	2.57	0.40
1:C:163:ILE:HG23	1:C:164:LEU:HD12	2.03	0.40
1:C:188:MET:C	1:C:190:LEU:N	2.78	0.40
1:D:202:GLN:O	1:D:203:PHE:CB	2.64	0.40
1:A:287:ASN:HA	1:A:318:LYS:HB2	2.03	0.40
1:B:190:LEU:HD22	1:B:193:MET:SD	2.61	0.40
1:C:182:GLY:HA2	1:C:204:ASP:HA	2.03	0.40
1:F:244:THR:HB	1:F:253:TYR:HB3	2.04	0.40
1:A:234:ILE:HD13	1:A:261:GLU:HG3	2.02	0.40
1:B:231:LYS:HA	1:B:231:LYS:HD3	1.91	0.40
1:C:135:VAL:HG22	4:C:432:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:VAL:HA	1:C:255:PRO:HD2	1.90	0.40
1:H:219:ASN:O	1:H:342:GLY:HA3	2.21	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:G:90:ASN:OD1	4:H:569:HOH:O[1_545]	2.19	0.01

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	309/347 (89%)	289 (94%)	18 (6%)	2 (1%)	21	32
1	B	309/347 (89%)	282 (91%)	27 (9%)	0	100	100
1	C	309/347 (89%)	287 (93%)	22 (7%)	0	100	100
1	D	309/347 (89%)	288 (93%)	20 (6%)	1 (0%)	36	50
1	E	309/347 (89%)	294 (95%)	15 (5%)	0	100	100
1	F	309/347 (89%)	283 (92%)	26 (8%)	0	100	100
1	G	309/347 (89%)	298 (96%)	11 (4%)	0	100	100
1	H	309/347 (89%)	279 (90%)	30 (10%)	0	100	100
All	All	2472/2776 (89%)	2300 (93%)	169 (7%)	3 (0%)	48	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	123	SER
1	A	144	SER
1	A	189	PRO

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	281/311 (90%)	266 (95%)	15 (5%)	20	36
1	B	281/311 (90%)	262 (93%)	19 (7%)	14	25
1	C	281/311 (90%)	262 (93%)	19 (7%)	14	25
1	D	281/311 (90%)	257 (92%)	24 (8%)	10	16
1	E	281/311 (90%)	268 (95%)	13 (5%)	24	41
1	F	281/311 (90%)	266 (95%)	15 (5%)	20	36
1	G	281/311 (90%)	265 (94%)	16 (6%)	18	33
1	H	281/311 (90%)	257 (92%)	24 (8%)	10	16
All	All	2248/2488 (90%)	2103 (94%)	145 (6%)	15	27

All (145) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	VAL
1	A	98	GLU
1	A	114	THR
1	A	119	ILE
1	A	136	THR
1	A	184	ILE
1	A	233	LEU
1	A	240	SER
1	A	285	THR
1	A	294	ASN
1	A	297	THR
1	A	303	LEU
1	A	305	VAL
1	A	328	SER
1	A	334	THR
1	B	48	ASP
1	B	98	GLU
1	B	104	LYS
1	B	114	THR

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Mol	Chain	Res	Type
1	B	125	LEU
1	B	146	LEU
1	B	151	LEU
1	B	189	PRO
1	B	233	LEU
1	B	240	SER
1	B	248	GLU
1	B	285	THR
1	B	294	ASN
1	B	297	THR
1	B	305	VAL
1	B	307	ASP
1	B	311	THR
1	B	325	ILE
1	B	328	SER
1	C	98	GLU
1	C	114	THR
1	C	125	LEU
1	C	140	ILE
1	C	146	LEU
1	C	151	LEU
1	C	169	THR
1	C	184	ILE
1	C	187	ILE
1	C	189	PRO
1	C	233	LEU
1	C	248	GLU
1	C	285	THR
1	C	294	ASN
1	C	297	THR
1	C	307	ASP
1	C	328	SER
1	C	347	SER
1	C	350	ILE
1	D	48	ASP
1	D	88	ASN
1	D	98	GLU
1	D	112	HIS
1	D	114	THR
1	D	140	ILE
1	D	146	LEU
1	D	164	LEU

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Mol	Chain	Res	Type
1	D	169	THR
1	D	184	ILE
1	D	201	ILE
1	D	233	LEU
1	D	285	THR
1	D	294	ASN
1	D	297	THR
1	D	299	THR
1	D	300	SER
1	D	303	LEU
1	D	305	VAL
1	D	307	ASP
1	D	315	ASN
1	D	334	THR
1	D	347	SER
1	D	357	VAL
1	E	98	GLU
1	E	125	LEU
1	E	176	ILE
1	E	233	LEU
1	E	240	SER
1	E	244	THR
1	E	285	THR
1	E	297	THR
1	E	303	LEU
1	E	305	VAL
1	E	307	ASP
1	E	347	SER
1	E	357	VAL
1	F	98	GLU
1	F	114	THR
1	F	119	ILE
1	F	187	ILE
1	F	191	LYS
1	F	233	LEU
1	F	240	SER
1	F	249	ASN
1	F	285	THR
1	F	294	ASN
1	F	315	ASN
1	F	328	SER
1	F	334	THR

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Mol	Chain	Res	Type
1	F	347	SER
1	F	357	VAL
1	G	56	THR
1	G	98	GLU
1	G	114	THR
1	G	129	ARG
1	G	146	LEU
1	G	167	ILE
1	G	169	THR
1	G	233	LEU
1	G	285	THR
1	G	294	ASN
1	G	297	THR
1	G	298	ILE
1	G	299	THR
1	G	307	ASP
1	G	334	THR
1	G	347	SER
1	H	69	SER
1	H	98	GLU
1	H	114	THR
1	H	136	THR
1	H	144	SER
1	H	146	LEU
1	H	151	LEU
1	H	167	ILE
1	H	169	THR
1	H	170	LEU
1	H	178	LEU
1	H	184	ILE
1	H	225	SER
1	H	233	LEU
1	H	240	SER
1	H	243	LEU
1	H	285	THR
1	H	294	ASN
1	H	297	THR
1	H	299	THR
1	H	307	ASP
1	H	315	ASN
1	H	328	SER
1	H	334	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	235	ASN
1	B	148	ASN
1	B	290	GLN
1	C	101	HIS
1	C	207	HIS
1	C	235	ASN
1	C	250	GLN
1	C	322	ASN
1	D	102	ASN
1	D	112	HIS
1	D	202	GLN
1	D	207	HIS
1	D	250	GLN
1	E	49	ASN
1	E	109	ASN
1	E	156	HIS
1	E	207	HIS
1	E	235	ASN
1	E	315	ASN
1	F	51	ASN
1	F	72	ASN
1	F	290	GLN
1	F	292	ASN
1	G	90	ASN
1	G	112	HIS
1	G	117	ASN
1	G	156	HIS
1	G	207	HIS
1	G	290	GLN
1	G	337	GLN
1	H	72	ASN
1	H	90	ASN
1	H	250	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	311/347 (89%)	-1.23	0 100 100	16, 30, 65, 82	0
1	B	311/347 (89%)	-1.28	0 100 100	13, 27, 58, 82	0
1	C	311/347 (89%)	-1.17	0 100 100	15, 29, 81, 99	0
1	D	311/347 (89%)	-1.12	0 100 100	17, 32, 79, 94	0
1	E	311/347 (89%)	-1.03	0 100 100	14, 39, 87, 97	0
1	F	311/347 (89%)	-1.25	0 100 100	14, 28, 67, 91	0
1	G	311/347 (89%)	-1.20	0 100 100	15, 33, 68, 87	0
1	H	311/347 (89%)	-1.27	0 100 100	12, 29, 66, 85	0
All	All	2488/2776 (89%)	-1.19	0 100 100	12, 31, 75, 99	0

There are no RSRZ outliers to report.

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](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	G	401	1/1	0.97	0.05	32,32,32,32	0
3	NA	H	401	1/1	0.98	0.06	29,29,29,29	0
2	NI	B	401	1/1	0.99	0.02	36,36,36,36	0
2	NI	F	401	1/1	1.00	0.02	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.