



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 06:12 AM GMT

PDB ID : 4L3F
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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

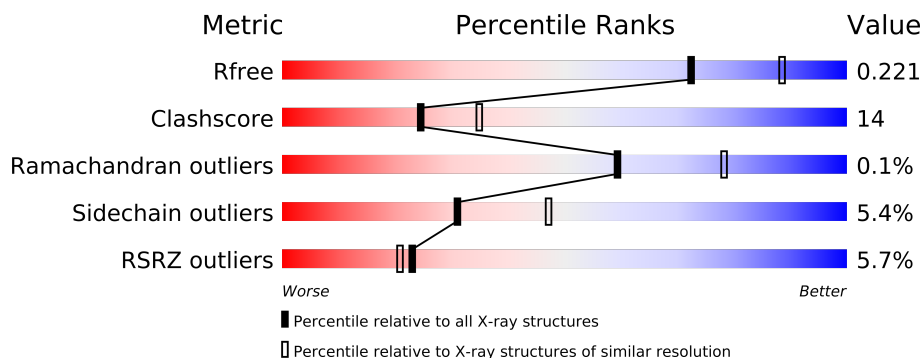
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	347	
1	B	347	
1	C	347	
1	D	347	
1	E	347	
1	F	347	
1	G	347	
1	H	347	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20171 atoms, of which 0 are hydrogen and 0 are deuterium.

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

Continued on next page...

Continued from previous page...

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

Continued on next page...

Continued from previous page...

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

Continued on next page...

Continued from previous page...

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

Continued on next page...

Continued from previous page...

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 (three-letter code: NI) (formula: Ni).

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
-----	-------	----------	-------	---------	---------

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 (three-letter code: NA) (formula: Na).

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

- Molecule 4 is water.

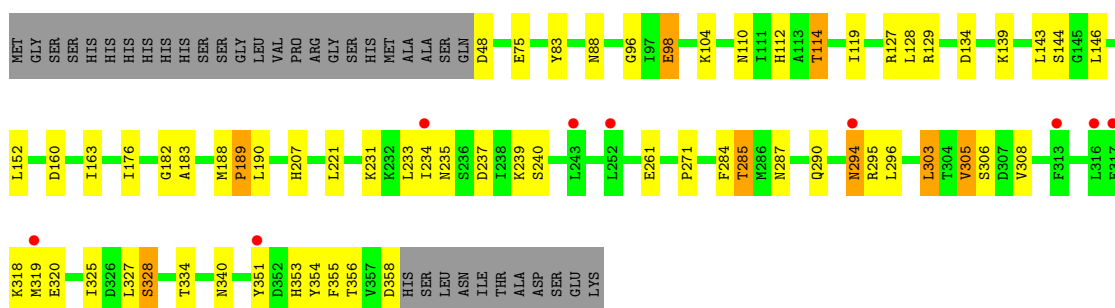
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	96	Total O 96 96	0	0
4	B	93	Total O 93 93	0	0
4	C	99	Total O 99 99	0	0
4	D	95	Total O 95 95	0	0
4	E	59	Total O 59 59	0	0
4	F	87	Total O 87 87	0	0
4	G	65	Total O 65 65	0	0
4	H	85	Total O 85 85	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

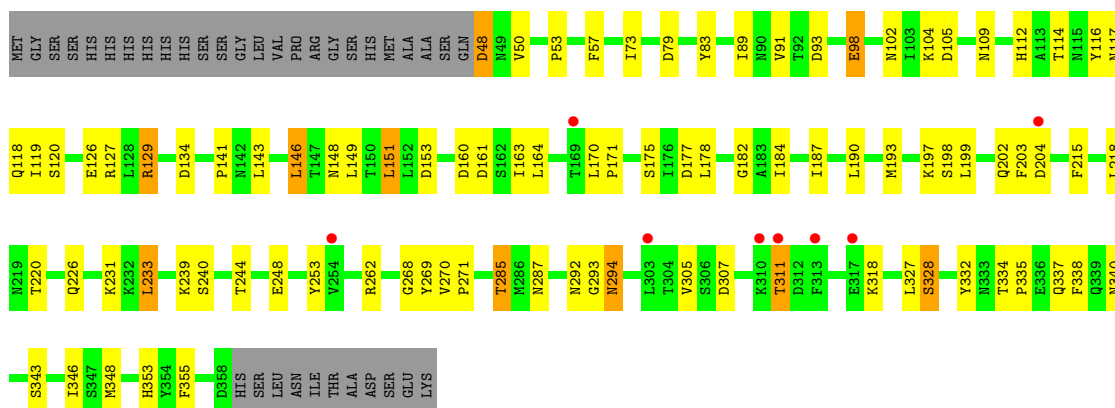
• Molecule 1: Internalin K

Chain A:



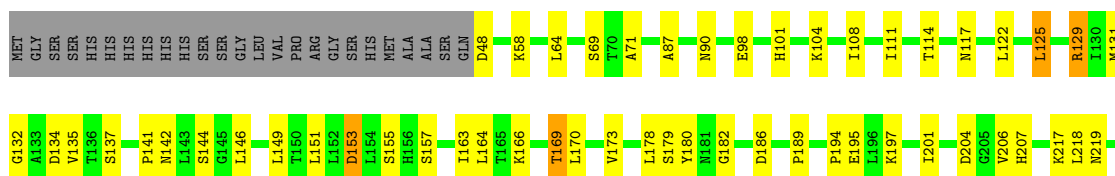
• Molecule 1: Internalin K

Chain B:

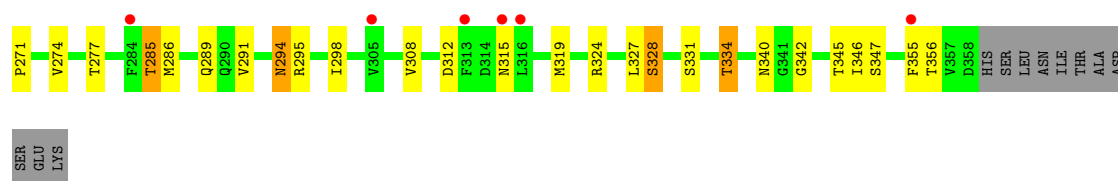


• Molecule 1: Internalin K

Chain C:

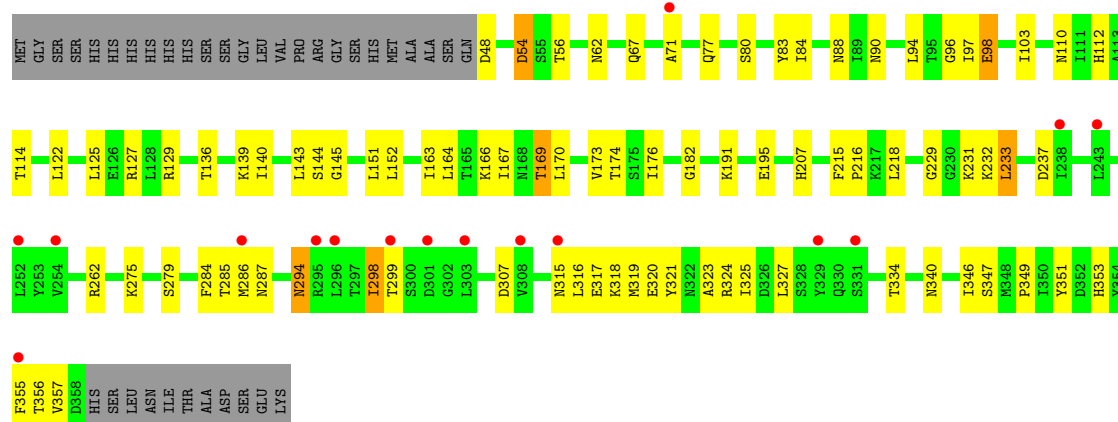






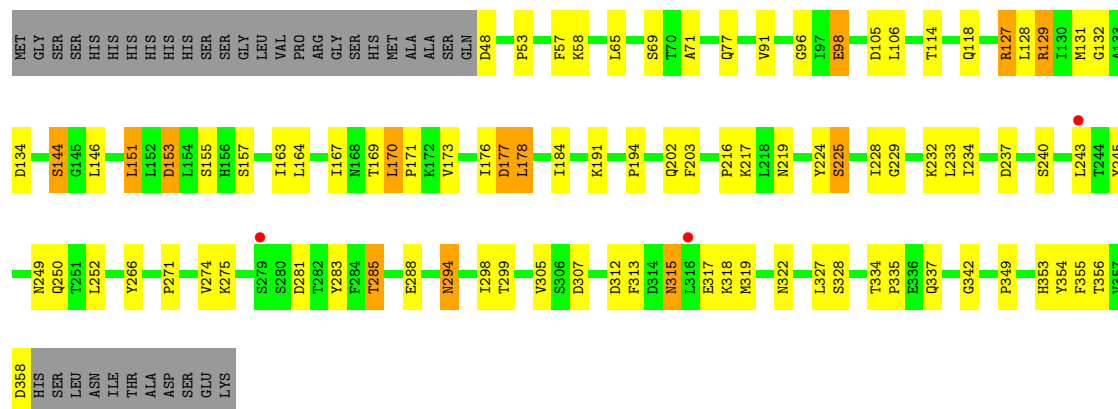
• Molecule 1: Internalin K

Chain G:



• Molecule 1: Internalin K

Chain H:



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 0.194 , 0.221	Depositor DCC
R_{free} test set	9215 reflections (11.10%)	DCC
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 , 11.8	EDS
Estimated twinning fraction	0.679 for H, K, L 0.321 for -H, -K, H+L 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
L-test for twinning	$\langle L \rangle = 0.35$, $\langle L^2 \rangle = 0.18$	Xtriage
Outliers	0 of 92214 reflections	Xtriage
F_o, F_c correlation	0.92	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.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.74	0/2478	0.87	1/3373 (0.0%)
1	B	0.83	0/2478	0.90	1/3373 (0.0%)
1	C	0.77	0/2478	0.89	4/3373 (0.1%)
1	D	0.74	0/2478	0.87	3/3373 (0.1%)
1	E	0.64	0/2478	0.80	1/3373 (0.0%)
1	F	0.79	0/2478	0.89	0/3373
1	G	0.74	1/2478 (0.0%)	0.90	1/3373 (0.0%)
1	H	0.78	0/2478	0.94	4/3373 (0.1%)
All	All	0.75	1/19824 (0.0%)	0.88	15/26984 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	83	TYR	CB-CG	5.25	1.59	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	127	ARG	NE-CZ-NH2	-7.70	116.45	120.30
1	D	177	ASP	CB-CG-OD1	6.44	124.10	118.30
1	H	129	ARG	NE-CZ-NH1	-6.18	117.21	120.30
1	C	129	ARG	NE-CZ-NH1	-5.89	117.35	120.30
1	A	127	ARG	NE-CZ-NH1	-5.85	117.37	120.30
1	H	153	ASP	CB-CG-OD1	5.70	123.43	118.30
1	C	151	LEU	CA-CB-CG	5.61	128.21	115.30
1	H	177	ASP	CB-CG-OD1	5.61	123.34	118.30
1	D	164	LEU	CB-CG-CD1	5.52	120.39	111.00
1	C	64	LEU	CB-CG-CD1	-5.49	101.66	111.00
1	G	54	ASP	CB-CG-OD2	5.30	123.07	118.30
1	E	177	ASP	CB-CG-OD1	5.26	123.03	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	129	ARG	CG-CD-NE	-5.25	100.78	111.80
1	D	64	LEU	CB-CG-CD1	-5.17	102.21	111.00
1	C	153	ASP	CB-CG-OD1	5.09	122.88	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2436	0	2375	48	0
1	B	2436	0	2375	81	0
1	C	2436	0	2375	82	0
1	D	2436	0	2375	90	0
1	E	2436	0	2375	62	0
1	F	2436	0	2375	59	0
1	G	2436	0	2375	62	1
1	H	2436	0	2375	66	0
2	B	1	0	0	0	0
2	F	1	0	0	0	0
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	27	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	537	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (537) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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: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:H:337:GLN:HG2	4:H:572:HOH:O	1.65	0.96
1:A:261:GLU:O	1:A:351:TYR:HE1	1.49	0.96
1:D:272:ASP:HA	4:D:451:HOH:O	1.65	0.96
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: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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:O	1:B:48:ASP:OD1	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: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
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:D:272:ASP:C	4:D:451:HOH:O	2.27	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:D:272:ASP:CA	4:D:451:HOH:O	2.30	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: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:D:275:LYS:HA	4:D:492:HOH:O	1.93	0.69
1:B:53:PRO:HD2	4:B:532:HOH:O	1.91	0.69
1:F:69:SER:O	4:F:520:HOH:O	2.08	0.69
1:C:90:ASN:HB2	4:C:481:HOH:O	1.93	0.69
1:D:202:GLN:O	1:D:203:PHE:HB2	1.93	0.69
1:C:329:TYR:CE1	4:C:477:HOH:O	2.42	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:339:GLN:O	4:E:401:HOH:O	2.11	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:G:340:ASN:HB2	4:G:557:HOH:O	1.94	0.68
1:E:120:SER:O	4:E:404:HOH:O	2.12	0.68
1:F:334:THR:HG23	4:F:568:HOH:O	1.93	0.67
1:F:200:ASN:OD1	1:F:202:GLN:HG3	1.94	0.67
1:B:204:ASP:O	1:B:226:GLN:OE1	2.13	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:D:262:ARG:NH2	1:D:271:PRO:O	2.28	0.66
1:A:285:THR:HG22	4:A:492:HOH:O	1.89	0.66
1:A:239:LYS:HD3	1:A:358:ASP:HB3	1.77	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:C:131:MET:CE	4:C:484:HOH:O	2.44	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
1:B:204:ASP:C	1:B:226:GLN:OE1	2.34	0.65
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:H:106:LEU:CB	4:H:584:HOH:O	2.44	0.65
1:C:271:PRO:HG3	1:C:327:LEU:HD13	1.77	0.65
1:H:294:ASN:N	1:H:294:ASN:OD1	2.27	0.65
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: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: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:G:340:ASN:CG	4:G:557:HOH:O	2.35	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
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: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:H:224:TYR:CD2	1:H:225:SER:HB2	2.36	0.60
1:G:122:LEU:O	1:G:125:LEU:HB2	2.01	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:H:202:GLN:O	1:H:203:PHE:HB2	2.02	0.60
1:B:262:ARG:NH2	1:B:271:PRO:O	2.33	0.60
1:C:246:ASN:OD1	1:C:249:ASN:HB2	2.02	0.60
1:F:183:ALA:HA	4:F:507:HOH:O	2.02	0.59
1:B:328:SER:HA	1:B:348:MET:CE	2.33	0.59
1:F:319:MET:HB2	1:F:355:PHE:HB2	1.83	0.59
1:B:120:SER:OG	1:C:142:ASN:CG	2.41	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:G:275:LYS:HD3	1:G:298:ILE:HG22	1.84	0.58
1:G:96:GLY:N	1:G:98:GLU:OE1	2.31	0.58
1:C:201:ILE:HD12	1:C:206:VAL:HG21	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: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:C:129:ARG:NH2	1:D:129:ARG:HE	2.02	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: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:G:77:GLN:O	1:G:80:SER:OG	2.16	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:319:MET:HB2	1:C:355:PHE:HB2	1.86	0.56
1:C:287:ASN:HA	1:C:318:LYS:HB2	1.88	0.56
1:C:197:LYS:O	1:C:218:LEU:HD12	2.06	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:G:287:ASN:ND2	1:G:315:ASN:O	2.39	0.55
1:E:104:LYS:NZ	4:E:426:HOH:O	2.37	0.55
1:G:317:GLU:O	1:G:356:THR:HA	2.06	0.55
1:G:77:GLN:HG2	4:G:527:HOH:O	2.07	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
1:F:277:THR:HG22	1:F:298:ILE:HD12	1.87	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:A:261:GLU:C	1:A:351:TYR:HE1	2.09	0.55
1:B:335:PRO:HB2	4:B:581:HOH:O	2.07	0.55
1:H:271:PRO:HG3	1:H:327:LEU:HD13	1.89	0.55
1:C:237:ASP:OD1	1:C:356:THR:OG1	2.14	0.54
1:H:48:ASP:N	4:H:507:HOH:O	2.39	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: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:B:50:VAL:HB	4:B:585:HOH:O	2.06	0.54
1:F:295:ARG:NH2	1:F:308:VAL:HG23	2.23	0.54
1:H:128:LEU:HD12	4:H:584:HOH:O	1.85	0.54
1:G:127:ARG:HG2	1:G:151:LEU:HB3	1.88	0.54
1:F:187:ILE:HD13	1:F:199:LEU:HD23	1.90	0.54
1:E:283:TYR:HB3	1:E:322:ASN:HB3	1.90	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:C:48:ASP:N	4:C:443:HOH:O	2.40	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:D:340:ASN:CG	4:D:474:HOH:O	2.46	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:B:328:SER:HA	1:B:348:MET:HE2	1.90	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:210:ARG:HA	4:F:534:HOH:O	2.08	0.53
1:F:340:ASN:O	1:F:340:ASN:CG	2.46	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
1:A:152:LEU:O	1:A:176:ILE:HA	2.09	0.53
1:A:114:THR:O	1:A:139:LYS:HE2	2.08	0.53
1:A:75:GLU:O	4:A:445:HOH:O	2.19	0.52
1:H:98:GLU:HG3	1:H:118:GLN:HB3	1.92	0.52
1:E:311:THR:HG22	1:E:315:ASN:ND2	2.24	0.52
1:D:246:ASN:OD1	1:D:249:ASN:N	2.28	0.52
1:D:122:LEU:C	1:D:124:ASN:N	2.63	0.52
1:A:261:GLU:C	1:A:351:TYR:CE1	2.83	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:E:319:MET:HB2	1:E:355:PHE:HB2	1.92	0.51
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:F:294:ASN:N	1:F:294:ASN:OD1	2.43	0.51
1:C:287:ASN:ND2	1:C:316:LEU:HD12	2.25	0.51
1:C:275:LYS:HD3	1:C:298:ILE:HG22	1.92	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:E:92:THR:HG22	1:E:92:THR:O	2.11	0.51
1:G:294:ASN:OD1	1:G:294:ASN:N	2.44	0.51
1:D:122:LEU:C	1:D:124:ASN:H	2.12	0.51
1:F:160:ASP:O	1:F:163:ILE:HG22	2.10	0.51
1:H:48:ASP:OD1	1:H:48:ASP:O	2.29	0.51
1:F:170:LEU:HB3	1:F:173:VAL:HG23	1.91	0.51
1:D:233:LEU:HB2	4:D:445:HOH:O	2.10	0.51
1:H:194:PRO:HD2	4:H:514:HOH:O	2.11	0.51
1:E:131:MET:HG2	1:E:155:SER:O	2.11	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:216:PRO:HD3	4:D:418:HOH:O	2.11	0.51
1:H:170:LEU:HG	1:H:173:VAL:HG21	1.92	0.51
1:E:237:ASP:OD1	1:E:356:THR:OG1	2.27	0.51
1:C:247:ALA:CB	4:C:482:HOH:O	2.58	0.51
1:C:224:TYR:OH	4:C:414:HOH:O	2.20	0.51
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:F:129:ARG:HH11	1:F:129:ARG:HG3	1.76	0.50
1:D:210:ARG:HD3	4:D:481:HOH:O	2.11	0.50
1:H:134:ASP:CG	4:H:520:HOH:O	2.49	0.50
1:D:286:MET:HG3	1:D:291:VAL:CG2	2.41	0.50
1:E:194:PRO:O	1:E:217:LYS:HD2	2.11	0.50
1:D:317:GLU:O	1:D:356:THR:HA	2.11	0.50
1:B:116:TYR:O	1:B:119:ILE:HG12	2.11	0.50
1:D:124:ASN:OD1	4:D:484:HOH:O	2.19	0.50
1:F:187:ILE:HG21	1:F:201:ILE:HG21	1.94	0.50
1:B:143:LEU:HB3	1:B:146:LEU:HD22	1.92	0.50
1:G:262:ARG:HB2	4:G:559:HOH:O	2.12	0.50
1:A:189:PRO:HD3	4:A:474:HOH:O	2.11	0.50
1:B:244:THR:HB	1:B:253:TYR:HB3	1.92	0.50
1:B:294:ASN:OD1	1:B:294:ASN:N	2.45	0.50
1:E:235:ASN:HA	1:E:354:TYR:O	2.12	0.50
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:D:117:ASN:HB3	4:D:462:HOH:O	2.12	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:H:128:LEU:HD13	4:H:584:HOH:O	1.91	0.49
1:D:286:MET:HA	4:D:494:HOH:O	2.12	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:B:79:ASP:O	1:B:102:ASN:ND2	2.43	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: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:C:87:ALA:HB3	1:D:88:ASN:HB2	1.95	0.49
1:A:295:ARG:O	1:A:305:VAL:HA	2.12	0.49
1:A:235:ASN:HA	1:A:354:TYR:O	2.12	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:E:143:LEU:HB3	1:E:146:LEU:HD22	1.95	0.48
1:H:202:GLN:HB3	1:H:203:PHE:CD2	2.48	0.48
1:A:182:GLY:HA3	1:A:231:LYS:HG3	1.95	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:G:140:ILE:O	1:G:166:LYS:NZ	2.39	0.48
1:H:129:ARG:HG2	1:H:153:ASP:HB3	1.94	0.48
1:C:207:HIS:O	1:C:263:THR:HA	2.13	0.48
1:E:271:PRO:HG3	1:E:327:LEU:HD13	1.95	0.48
1:D:170:LEU:HB3	1:D:173:VAL:HG23	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: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:B:178:LEU:HD13	1:B:184:ILE:HD13	1.95	0.48
1:C:129:ARG:NH2	1:D:129:ARG:NH2	2.62	0.47
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:H:144:SER:O	4:H:558:HOH:O	2.20	0.47
1:D:62:ASN:OD1	1:D:67:GLN:NE2	2.47	0.47
1:H:176:ILE:HG22	1:H:177:ASP:N	2.29	0.47
1:D:339:GLN:O	4:D:403:HOH:O	2.20	0.47
1:H:245:TYR:HB2	1:H:313:PHE:CE1	2.49	0.47
1:G:353:HIS:CD2	1:G:355:PHE:CZ	3.03	0.47
1:E:296:LEU:HD23	1:E:305:VAL:HG13	1.97	0.47
1:A:188:MET:O	1:A:190:LEU:N	2.47	0.47
1:F:58:LYS:HE2	1:F:71:ALA:O	2.14	0.47
1:H:96:GLY:N	1:H:98:GLU:OE1	2.43	0.47
1:A:284:PHE:O	1:A:290:GLN:HA	2.15	0.47
1:A:143:LEU:HB3	1:A:146:LEU:HD22	1.97	0.47
1:D:48:ASP:O	1:D:48:ASP:OD1	2.33	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:F:143:LEU:HD12	1:F:167:ILE:HG22	1.95	0.47
1:B:105:ASP:OD1	1:B:127:ARG:CZ	2.63	0.47
1:D:244:THR:HB	1:D:253:TYR:HB3	1.97	0.47
1:E:236:SER:HB2	1:E:355:PHE:HD1	1.77	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:203:PHE:HA	1:H:225:SER:O	2.14	0.47
1:B:129:ARG:HG2	1:B:153:ASP:HB3	1.97	0.47
1:H:237:ASP:OD1	1:H:356:THR:OG1	2.26	0.47
1:H:315:ASN:OD1	1:H:315:ASN:N	2.47	0.47
1:G:94:LEU:O	1:G:97:ILE:HG22	2.13	0.47
1:C:129:ARG:NH2	1:D:129:ARG:CZ	2.77	0.46
1:E:105:ASP:OD1	1:E:127:ARG:HB2	2.15	0.46
1:E:125:LEU:HA	1:E:125:LEU:HD12	1.67	0.46
1:E:344:TYR:N	4:E:456:HOH:O	2.47	0.46
1:H:219:ASN:C	1:H:219:ASN:OD1	2.54	0.46
1:C:322:ASN:HA	1:C:351:TYR:O	2.15	0.46
1:D:253:TYR:CE1	1:D:302:GLY:HA3	2.51	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:G:174:THR:HA	1:G:195:GLU:O	2.16	0.46
1:C:134:ASP:OD1	1:C:134:ASP:N	2.49	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:H:266:TYR:CD2	1:H:335:PRO:HG3	2.50	0.46
1:F:103:ILE:HG23	4:F:506:HOH:O	2.16	0.46
1:B:202:GLN:O	1:B:203:PHE:HB2	2.15	0.46
1:C:251:THR:HA	1:C:305:VAL:O	2.15	0.46
1:E:152:LEU:HD23	1:E:176:ILE:HG12	1.97	0.46
1:A:112:HIS:HA	1:A:134:ASP:OD2	2.16	0.46
1:D:274:VAL:HG21	1:D:281:ASP:HB3	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:E:52:ILE:HB	1:E:58:LYS:HD2	1.98	0.45
1:D:92:THR:HG22	1:D:92:THR:O	2.16	0.45
1:B:311:THR:HG23	4:B:540:HOH:O	2.16	0.45
1:G:110:ASN:OD1	1:G:112:HIS:HE1	1.98	0.45
1:B:164:LEU:HD23	1:B:190:LEU:HG	1.97	0.45
1:B:126:GLU:O	1:B:149:LEU:HD12	2.17	0.45
1:H:129:ARG:HG3	1:H:129:ARG:HH11	1.80	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:D:166:LYS:HG2	4:D:463:HOH:O	2.16	0.45
1:C:294:ASN:OD1	1:C:294:ASN:N	2.49	0.45
1:D:251:THR:HA	1:D:305:VAL:O	2.16	0.45
1:B:353:HIS:CD2	1:B:355:PHE:CZ	3.05	0.45
1:B:268:GLY:O	1:B:270:VAL:HG23	2.16	0.45
1:F:295:ARG:HH22	1:F:308:VAL:CG2	2.29	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:H:57:PHE:HB2	1:H:91:VAL:HG13	1.98	0.45
1:B:182:GLY:HA3	1:B:231:LYS:HG2	1.99	0.45
1:C:186:ASP:C	1:C:186:ASP:OD1	2.54	0.45
1:B:340:ASN:CG	4:B:543:HOH:O	2.45	0.45
1:F:219:ASN:O	1:F:342:GLY:HA3	2.17	0.45
1:C:178:LEU:HA	1:C:178:LEU:HD23	1.82	0.45
1:G:125:LEU:HA	1:G:125:LEU:HD12	1.77	0.45
1:D:153:ASP:C	1:D:153:ASP:OD1	2.54	0.45
1:D:275:LYS:CA	4:D:492:HOH:O	2.60	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:D:204:ASP:O	1:D:226:GLN:HA	2.17	0.44
1:H:234:ILE:HG21	1:H:353:HIS:CE1	2.52	0.44
1:B:187:ILE:HD12	1:B:199:LEU:CD2	2.47	0.44
1:B:127:ARG:HG2	1:B:151:LEU:HB3	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:C:108:ILE:HG23	1:C:111:ILE:HG21	1.99	0.44
1:A:96:GLY:CA	1:A:98:GLU:OE1	2.65	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:F:223:ALA:HB3	1:F:346:ILE:HG22	2.00	0.44
1:B:292:ASN:OD1	1:B:293:GLY:N	2.50	0.44
1:C:149:LEU:HA	1:C:149:LEU:HD12	1.77	0.44
1:G:191:LYS:HA	1:G:215:PHE:CE1	2.52	0.44
1:B:120:SER:HB3	1:B:141:PRO:HG2	1.99	0.44
1:B:328:SER:HA	1:B:348:MET:HE1	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:319:MET:HB2	1:G:355:PHE:HB2	1.98	0.44
1:G:173:VAL:HG11	1:G:176:ILE:HG13	1.99	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:319:MET:HB2	1:D:355:PHE:HB2	1.99	0.44
1:D:245:TYR:OH	1:D:308:VAL:O	2.24	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:C:170:LEU:HD13	1:C:173:VAL:HG21	2.00	0.44
1:H:285:THR:O	1:H:319:MET:HA	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:271:PRO:HB3	1:D:327:LEU:HD13	2.00	0.43
1:C:285:THR:HB	1:C:320:GLU:HB3	1.99	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:C:117:ASN:HA	1:C:141:PRO:HG3	2.00	0.43
1:H:170:LEU:HA	1:H:171:PRO:HD3	1.75	0.43
1:D:100:ALA:C	4:D:480:HOH:O	2.56	0.43
1:A:83:TYR:HE1	4:A:457:HOH:O	2.02	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:C:328:SER:CB	4:C:486:HOH:O	2.07	0.43
1:G:151:LEU:C	1:G:151:LEU:HD23	2.39	0.43
1:G:62:ASN:OD1	1:G:71:ALA:O	2.36	0.43
1:B:146:LEU:HD23	1:B:149:LEU:HD22	2.01	0.43
1:F:286:MET:HG3	1:F:291:VAL:CG2	2.48	0.43
1:F:274:VAL:HG22	1:F:324:ARG:O	2.19	0.43
1:E:191:LYS:HB3	1:E:191:LYS:HE2	1.79	0.43
1:H:129:ARG:NH1	1:H:129:ARG:HG3	2.33	0.43
1:E:122:LEU:O	1:E:125:LEU:HB2	2.19	0.43
1:H:353:HIS:CD2	1:H:355:PHE:CZ	3.06	0.43
1:G:170:LEU:HD13	1:G:173:VAL:HG21	2.00	0.43
1:D:292:ASN:HB3	1:D:295:ARG:HG3	2.00	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:C:69:SER:O	4:C:447:HOH:O	2.20	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:C:131:MET:HE2	4:C:484:HOH:O	2.15	0.42
1:G:287:ASN:HD21	1:G:316:LEU:HA	1.82	0.42
1:H:317:GLU:O	1:H:356:THR:HA	2.19	0.42
1:C:219:ASN:C	1:C:219:ASN:OD1	2.57	0.42
1:B:287:ASN:HA	1:B:318:LYS:HB2	2.00	0.42
1:C:129:ARG:HH22	1:D:129:ARG:HH21	1.68	0.42
1:H:202:GLN:O	1:H:203:PHE:CB	2.67	0.42
1:D:237:ASP:OD1	1:D:356:THR:OG1	2.33	0.42
1:A:319:MET:HB2	1:A:355:PHE:HB2	2.01	0.42
1:G:316:LEU:HD23	1:G:357:VAL:HG21	2.01	0.42
1:D:129:ARG:HG3	1:D:129:ARG:HH11	1.84	0.42
1:F:334:THR:CG2	4:F:568:HOH:O	2.60	0.42
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:207:HIS:HD2	4:A:448:HOH:O	2.01	0.42
1:D:286:MET:O	1:D:289:GLN:HB2	2.18	0.42
1:F:197:LYS:C	1:F:218:LEU:HD12	2.40	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: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:C:229:GLY:HA2	1:C:232:LYS:O	2.20	0.42
1:B:170:LEU:HA	1:B:171:PRO:HD2	1.95	0.42
1:E:294:ASN:OD1	1:E:294:ASN:N	2.52	0.42
1:G:144:SER:HB3	1:G:169:THR:OG1	2.19	0.42
1:G:48:ASP:OD1	1:G:48:ASP:O	2.37	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:H:275:LYS:HD3	1:H:298:ILE:HG22	2.00	0.42
1:G:152:LEU:HD12	1:G:152:LEU:HA	1.81	0.42
1:B:149:LEU:HA	1:B:149:LEU:HD12	1.81	0.42
1:D:176:ILE:HD11	1:D:193:MET:CE	2.50	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:A:221:LEU:HA	1:A:221:LEU:HD12	1.90	0.42
1:G:315:ASN:CB	4:G:556:HOH:O	2.67	0.42
1:D:210:ARG:CD	4:D:481:HOH:O	2.67	0.42
1:H:53:PRO:HG3	4:H:546:HOH:O	2.20	0.42
1:C:286:MET:HB2	1:C:286:MET:HE2	1.95	0.42
1:D:94:LEU:HD23	1:D:94:LEU:HA	1.79	0.42
1:E:236:SER:HB2	1:E:355:PHE:CE1	2.54	0.41
1:H:327:LEU:HB3	1:H:349:PRO:HD2	2.02	0.41
1:H:178:LEU:HA	1:H:178:LEU:HD12	1.97	0.41
1:F:99:TYR:HA	1:F:101:HIS:CE1	2.55	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:E:120:SER:HA	1:E:146:LEU:HD11	2.02	0.41
1:A:340:ASN:ND2	4:A:414:HOH:O	2.40	0.41
1:E:268:GLY:O	1:E:270:VAL:HG23	2.20	0.41
1:D:204:ASP:C	1:D:226:GLN:OE1	2.58	0.41
1:C:129:ARG:HH21	1:D:129:ARG:CZ	2.33	0.41
1:G:191:LYS:HA	1:G:215:PHE:CD1	2.56	0.41
1:C:170:LEU:HB3	1:C:173:VAL:HG23	2.01	0.41
1:G:94:LEU:HA	1:G:94:LEU:HD23	1.81	0.41
1:E:74:THR:CG2	4:E:435:HOH:O	2.68	0.41
1:H:274:VAL:HG21	1:H:281:ASP:HB3	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:271:PRO:CB	1:F:327:LEU:HD13	2.50	0.41
1:B:175:SER:CB	4:B:572:HOH:O	2.59	0.41
1:C:122:LEU:CB	1:C:125:LEU:HD22	2.47	0.41
1:B:197:LYS:HA	1:B:218:LEU:HA	2.02	0.41
1:E:164:LEU:HD21	1:E:187:ILE:HA	2.03	0.41
1:E:239:LYS:HA	1:E:239:LYS:HD3	1.88	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: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:D:170:LEU:HA	1:D:170:LEU:HD23	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:G:163:ILE:HD12	1:G:163:ILE:HA	1.93	0.40
1:H:176:ILE:CG2	1:H:177:ASP:N	2.84	0.40
1:D:212:ILE:O	1:D:215:PHE:HB2	2.20	0.40
1:G:218:LEU:N	4:G:563:HOH:O	2.49	0.40
1:A:325:ILE:HD11	1:A:351:TYR:HB2	2.02	0.40
1:D:202:GLN:O	1:D:203:PHE:CB	2.64	0.40
1:A:296:LEU:HD22	1:A:303:LEU:HD11	2.03	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:B:233:LEU:HD12	1:B:233:LEU:HA	1.84	0.40
1:B:190:LEU:HD22	1:B:193:MET:SD	2.61	0.40
1:A:287:ASN:HA	1:A:318:LYS:HB2	2.03	0.40
1:F:244:THR:HB	1:F:253:TYR:HB3	2.04	0.40
1:C:182:GLY:HA2	1:C:204:ASP:HA	2.03	0.40
1:A:234:ILE:HD13	1:A:261:GLU:HG3	2.02	0.40
1:H:219:ASN:O	1:H:342:GLY:HA3	2.21	0.40
1:B:231:LYS:HD3	1:B:231:LYS:HA	1.91	0.40
1:C:135:VAL:HG22	4:C:432:HOH:O	2.21	0.40
1:C:254:VAL:HA	1:C:255:PRO:HD2	1.90	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	Distance(Å)	Clash(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%)	33	47
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%)	50	68
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%)	59	78

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%)	271 (96%)	10 (4%)	47	68
1	B	281/311 (90%)	265 (94%)	16 (6%)	29	44
1	C	281/311 (90%)	268 (95%)	13 (5%)	37	55
1	D	281/311 (90%)	263 (94%)	18 (6%)	25	37

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	281/311 (90%)	269 (96%)	12 (4%)	40	59
1	F	281/311 (90%)	268 (95%)	13 (5%)	37	55
1	G	281/311 (90%)	265 (94%)	16 (6%)	29	44
1	H	281/311 (90%)	258 (92%)	23 (8%)	17	24
All	All	2248/2488 (90%)	2127 (95%)	121 (5%)	31	47

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	GLU
1	A	114	THR
1	A	119	ILE
1	A	233	LEU
1	A	240	SER
1	A	285	THR
1	A	294	ASN
1	A	303	LEU
1	A	305	VAL
1	A	328	SER
1	B	48	ASP
1	B	57	PHE
1	B	98	GLU
1	B	104	LYS
1	B	114	THR
1	B	146	LEU
1	B	151	LEU
1	B	233	LEU
1	B	240	SER
1	B	248	GLU
1	B	285	THR
1	B	294	ASN
1	B	305	VAL
1	B	307	ASP
1	B	311	THR
1	B	328	SER
1	C	98	GLU
1	C	114	THR
1	C	125	LEU
1	C	146	LEU
1	C	169	THR
1	C	189	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	233	LEU
1	C	285	THR
1	C	294	ASN
1	C	303	LEU
1	C	307	ASP
1	C	328	SER
1	C	347	SER
1	D	48	ASP
1	D	98	GLU
1	D	112	HIS
1	D	114	THR
1	D	146	LEU
1	D	164	LEU
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	299	THR
1	D	300	SER
1	D	303	LEU
1	D	315	ASN
1	D	334	THR
1	D	347	SER
1	E	53	PRO
1	E	98	GLU
1	E	125	LEU
1	E	233	LEU
1	E	240	SER
1	E	285	THR
1	E	294	ASN
1	E	303	LEU
1	E	305	VAL
1	E	307	ASP
1	E	347	SER
1	E	357	VAL
1	F	114	THR
1	F	119	ILE
1	F	187	ILE
1	F	191	LYS
1	F	233	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
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
1	F	347	SER
1	G	56	THR
1	G	88	ASN
1	G	98	GLU
1	G	114	THR
1	G	129	ARG
1	G	167	ILE
1	G	169	THR
1	G	233	LEU
1	G	237	ASP
1	G	279	SER
1	G	294	ASN
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	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	191	LYS
1	H	225	SER
1	H	233	LEU
1	H	240	SER
1	H	243	LEU
1	H	285	THR
1	H	294	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	299	THR
1	H	307	ASP
1	H	315	ASN
1	H	328	SER
1	H	358	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	148	ASN
1	C	101	HIS
1	C	207	HIS
1	C	250	GLN
1	D	112	HIS
1	D	156	HIS
1	D	202	GLN
1	D	250	GLN
1	E	156	HIS
1	E	207	HIS
1	E	315	ASN
1	F	51	ASN
1	F	72	ASN
1	F	292	ASN
1	F	322	ASN
1	G	88	ASN
1	G	90	ASN
1	G	112	HIS
1	G	117	ASN
1	G	207	HIS
1	H	72	ASN
1	H	90	ASN
1	H	250	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	311/347 (89%)	0.38	9 (2%) 49 47	16, 30, 65, 82	0
1	B	311/347 (89%)	0.37	8 (2%) 53 51	13, 27, 58, 83	0
1	C	311/347 (89%)	0.52	17 (5%) 24 22	15, 29, 81, 99	0
1	D	311/347 (89%)	0.66	33 (10%) 7 6	17, 32, 79, 94	0
1	E	311/347 (89%)	0.88	46 (14%) 3 2	14, 39, 87, 97	0
1	F	311/347 (89%)	0.32	10 (3%) 45 43	14, 28, 67, 91	0
1	G	311/347 (89%)	0.49	16 (5%) 27 25	15, 33, 68, 87	0
1	H	311/347 (89%)	0.28	3 (0%) 79 79	12, 29, 66, 85	0
All	All	2488/2776 (89%)	0.49	142 (5%) 23 21	12, 31, 75, 99	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	316	LEU	7.4
1	C	243	LEU	6.2
1	E	304	THR	5.6
1	C	329	TYR	5.2
1	G	329	TYR	5.1
1	D	252	LEU	5.1
1	E	284	PHE	5.0
1	D	300	SER	4.7
1	C	244	THR	4.7
1	D	238	ILE	4.6
1	D	293	GLY	4.5
1	E	243	LEU	4.4
1	A	316	LEU	4.4
1	E	357	VAL	4.4
1	E	355	PHE	4.4
1	D	306	SER	4.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	294	ASN	4.2
1	E	252	LEU	4.1
1	E	317	GLU	4.0
1	C	313	PHE	3.9
1	C	284	PHE	3.9
1	E	245	TYR	3.8
1	G	254	VAL	3.7
1	E	291	VAL	3.7
1	D	241	SER	3.6
1	D	329	TYR	3.5
1	D	297	THR	3.5
1	E	247	ALA	3.5
1	G	296	LEU	3.4
1	D	305	VAL	3.4
1	D	286	MET	3.3
1	E	286	MET	3.3
1	E	254	VAL	3.3
1	D	304	THR	3.3
1	E	319	MET	3.3
1	E	306	SER	3.3
1	E	274	VAL	3.2
1	C	252	LEU	3.2
1	E	329	TYR	3.2
1	H	243	LEU	3.2
1	A	313	PHE	3.2
1	A	243	LEU	3.1
1	D	311	THR	3.1
1	C	253	TYR	3.1
1	G	308	VAL	3.1
1	G	252	LEU	3.0
1	C	249	ASN	3.0
1	G	243	LEU	3.0
1	D	283	TYR	3.0
1	D	243	LEU	3.0
1	D	203	PHE	2.9
1	E	244	THR	2.9
1	E	279	SER	2.9
1	E	300	SER	2.9
1	E	236	SER	2.9
1	F	316	LEU	2.9
1	C	241	SER	2.9
1	A	252	LEU	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	351	TYR	2.8
1	F	239	LYS	2.8
1	C	357	VAL	2.8
1	D	111	ILE	2.8
1	E	313	PHE	2.8
1	G	331	SER	2.7
1	E	305	VAL	2.7
1	G	238	ILE	2.7
1	E	237	ASP	2.7
1	B	310	LYS	2.7
1	F	315	ASN	2.7
1	E	349	PRO	2.6
1	F	284	PHE	2.6
1	D	303	LEU	2.6
1	G	301	ASP	2.6
1	G	286	MET	2.6
1	E	282	THR	2.6
1	B	303	LEU	2.6
1	D	249	ASN	2.6
1	E	238	ILE	2.5
1	B	313	PHE	2.5
1	D	323	ALA	2.5
1	E	253	TYR	2.5
1	E	314	ASP	2.5
1	C	237	ASP	2.5
1	D	307	ASP	2.5
1	A	294	ASN	2.5
1	E	275	LYS	2.5
1	E	298	ILE	2.5
1	F	355	PHE	2.4
1	H	316	LEU	2.4
1	D	124	ASN	2.4
1	E	315	ASN	2.4
1	D	319	MET	2.4
1	E	318	LYS	2.4
1	E	182	GLY	2.4
1	E	248	GLU	2.4
1	E	320	GLU	2.4
1	D	316	LEU	2.4
1	G	295	ARG	2.4
1	D	254	VAL	2.4
1	G	355	PHE	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	247	ALA	2.3
1	E	276	SER	2.3
1	C	291	VAL	2.3
1	C	282	THR	2.3
1	C	319	MET	2.3
1	E	277	THR	2.3
1	C	293	GLY	2.3
1	G	71	ALA	2.3
1	E	294	ASN	2.3
1	B	311	THR	2.3
1	C	283	TYR	2.3
1	G	299	THR	2.3
1	C	238	ILE	2.3
1	F	238	ILE	2.3
1	D	284	PHE	2.3
1	A	319	MET	2.2
1	G	303	LEU	2.2
1	E	350	ILE	2.2
1	E	321	TYR	2.2
1	D	336	GLU	2.2
1	F	248	GLU	2.2
1	F	313	PHE	2.2
1	D	324	ARG	2.1
1	B	204	ASP	2.1
1	D	296	LEU	2.1
1	F	258	LEU	2.1
1	G	315	ASN	2.1
1	A	317	GLU	2.1
1	B	254	VAL	2.1
1	D	298	ILE	2.1
1	B	169	THR	2.1
1	B	317	GLU	2.1
1	D	278	ALA	2.1
1	F	305	VAL	2.1
1	A	351	TYR	2.1
1	E	250	GLN	2.0
1	E	251	THR	2.0
1	H	279	SER	2.0
1	E	278	ALA	2.0
1	E	185	THR	2.0
1	A	234	ILE	2.0
1	D	239	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NA	H	401	1/1	0.13	-1.55	29,29,29,29	0
2	NI	B	401	1/1	0.11	-2.06	36,36,36,36	0
2	NI	F	401	1/1	0.06	-3.38	40,40,40,40	0
3	NA	G	401	1/1	0.10	-4.26	32,32,32,32	0

6.5 Other polymers ⓘ

There are no such residues in this entry.