



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:49 PM GMT

PDB ID : 3P5T
Title : CFIm25-CFIm68 complex
Authors : Li, H.; Tong, S.; Li, X.; Shi, H.; Gao, Y.; Ge, H.; Niu, L.; Teng, M.
Deposited on : 2010-10-11
Resolution : 2.70 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>

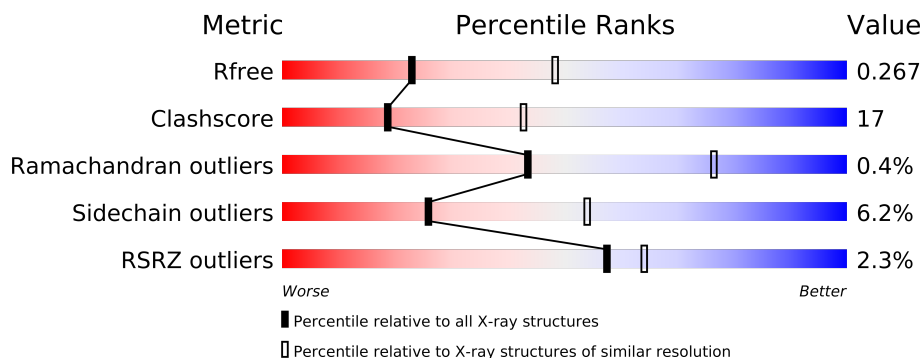
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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	202	
1	B	202	
1	C	202	
1	D	202	
1	E	202	
1	F	202	
2	L	90	
2	M	90	
2	N	90	
2	O	90	
2	P	90	
2	Q	90	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13253 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 Cleavage and polyadenylation specificity factor subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1589	1037	270	278	4			
1	B	198	Total	C	N	O	S	0	0	0
			1611	1048	273	286	4			
1	C	196	Total	C	N	O	S	0	0	0
			1574	1028	264	278	4			
1	D	198	Total	C	N	O	S	0	0	0
			1613	1052	272	285	4			
1	E	191	Total	C	N	O	S	0	0	0
			1519	994	253	268	4			
1	F	197	Total	C	N	O	S	0	0	0
			1597	1041	270	282	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	LEU	-	EXPRESSION TAG	UNP O43809
A	229	GLU	-	EXPRESSION TAG	UNP O43809
A	230	HIS	-	EXPRESSION TAG	UNP O43809
A	231	HIS	-	EXPRESSION TAG	UNP O43809
A	232	HIS	-	EXPRESSION TAG	UNP O43809
A	233	HIS	-	EXPRESSION TAG	UNP O43809
A	234	HIS	-	EXPRESSION TAG	UNP O43809
A	235	HIS	-	EXPRESSION TAG	UNP O43809
B	228	LEU	-	EXPRESSION TAG	UNP O43809
B	229	GLU	-	EXPRESSION TAG	UNP O43809
B	230	HIS	-	EXPRESSION TAG	UNP O43809
B	231	HIS	-	EXPRESSION TAG	UNP O43809
B	232	HIS	-	EXPRESSION TAG	UNP O43809
B	233	HIS	-	EXPRESSION TAG	UNP O43809
B	234	HIS	-	EXPRESSION TAG	UNP O43809
B	235	HIS	-	EXPRESSION TAG	UNP O43809
C	228	LEU	-	EXPRESSION TAG	UNP O43809

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Chain	Residue	Modelled	Actual	Comment	Reference
C	229	GLU	-	EXPRESSION TAG	UNP O43809
C	230	HIS	-	EXPRESSION TAG	UNP O43809
C	231	HIS	-	EXPRESSION TAG	UNP O43809
C	232	HIS	-	EXPRESSION TAG	UNP O43809
C	233	HIS	-	EXPRESSION TAG	UNP O43809
C	234	HIS	-	EXPRESSION TAG	UNP O43809
C	235	HIS	-	EXPRESSION TAG	UNP O43809
D	228	LEU	-	EXPRESSION TAG	UNP O43809
D	229	GLU	-	EXPRESSION TAG	UNP O43809
D	230	HIS	-	EXPRESSION TAG	UNP O43809
D	231	HIS	-	EXPRESSION TAG	UNP O43809
D	232	HIS	-	EXPRESSION TAG	UNP O43809
D	233	HIS	-	EXPRESSION TAG	UNP O43809
D	234	HIS	-	EXPRESSION TAG	UNP O43809
D	235	HIS	-	EXPRESSION TAG	UNP O43809
E	228	LEU	-	EXPRESSION TAG	UNP O43809
E	229	GLU	-	EXPRESSION TAG	UNP O43809
E	230	HIS	-	EXPRESSION TAG	UNP O43809
E	231	HIS	-	EXPRESSION TAG	UNP O43809
E	232	HIS	-	EXPRESSION TAG	UNP O43809
E	233	HIS	-	EXPRESSION TAG	UNP O43809
E	234	HIS	-	EXPRESSION TAG	UNP O43809
E	235	HIS	-	EXPRESSION TAG	UNP O43809
F	228	LEU	-	EXPRESSION TAG	UNP O43809
F	229	GLU	-	EXPRESSION TAG	UNP O43809
F	230	HIS	-	EXPRESSION TAG	UNP O43809
F	231	HIS	-	EXPRESSION TAG	UNP O43809
F	232	HIS	-	EXPRESSION TAG	UNP O43809
F	233	HIS	-	EXPRESSION TAG	UNP O43809
F	234	HIS	-	EXPRESSION TAG	UNP O43809
F	235	HIS	-	EXPRESSION TAG	UNP O43809

- Molecule 2 is a protein called Cleavage and polyadenylation specificity factor subunit 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	76	Total	C	N	O	S	0	0	0
			527	337	92	97	1			
2	M	81	Total	C	N	O	S	0	0	0
			603	385	100	117	1			
2	N	80	Total	C	N	O	S	0	0	0
			604	386	103	114	1			
2	O	81	Total	C	N	O	S	0	0	0
			599	382	98	118	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	79	Total	C	N	O	S	0	0	0
			586	375	97	113	1			
2	Q	80	Total	C	N	O	S	0	0	0
			583	372	97	113	1			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
L	162	LEU	-	EXPRESSION TAG	UNP Q16630
L	163	GLU	-	EXPRESSION TAG	UNP Q16630
L	164	HIS	-	EXPRESSION TAG	UNP Q16630
L	165	HIS	-	EXPRESSION TAG	UNP Q16630
L	166	HIS	-	EXPRESSION TAG	UNP Q16630
L	167	HIS	-	EXPRESSION TAG	UNP Q16630
L	168	HIS	-	EXPRESSION TAG	UNP Q16630
L	169	HIS	-	EXPRESSION TAG	UNP Q16630
M	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
M	162	LEU	-	EXPRESSION TAG	UNP Q16630
M	163	GLU	-	EXPRESSION TAG	UNP Q16630
M	164	HIS	-	EXPRESSION TAG	UNP Q16630
M	165	HIS	-	EXPRESSION TAG	UNP Q16630
M	166	HIS	-	EXPRESSION TAG	UNP Q16630
M	167	HIS	-	EXPRESSION TAG	UNP Q16630
M	168	HIS	-	EXPRESSION TAG	UNP Q16630
M	169	HIS	-	EXPRESSION TAG	UNP Q16630
N	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
N	162	LEU	-	EXPRESSION TAG	UNP Q16630
N	163	GLU	-	EXPRESSION TAG	UNP Q16630
N	164	HIS	-	EXPRESSION TAG	UNP Q16630
N	165	HIS	-	EXPRESSION TAG	UNP Q16630
N	166	HIS	-	EXPRESSION TAG	UNP Q16630
N	167	HIS	-	EXPRESSION TAG	UNP Q16630
N	168	HIS	-	EXPRESSION TAG	UNP Q16630
N	169	HIS	-	EXPRESSION TAG	UNP Q16630
O	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
O	162	LEU	-	EXPRESSION TAG	UNP Q16630
O	163	GLU	-	EXPRESSION TAG	UNP Q16630
O	164	HIS	-	EXPRESSION TAG	UNP Q16630
O	165	HIS	-	EXPRESSION TAG	UNP Q16630
O	166	HIS	-	EXPRESSION TAG	UNP Q16630
O	167	HIS	-	EXPRESSION TAG	UNP Q16630

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Chain	Residue	Modelled	Actual	Comment	Reference
O	168	HIS	-	EXPRESSION TAG	UNP Q16630
O	169	HIS	-	EXPRESSION TAG	UNP Q16630
P	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
P	162	LEU	-	EXPRESSION TAG	UNP Q16630
P	163	GLU	-	EXPRESSION TAG	UNP Q16630
P	164	HIS	-	EXPRESSION TAG	UNP Q16630
P	165	HIS	-	EXPRESSION TAG	UNP Q16630
P	166	HIS	-	EXPRESSION TAG	UNP Q16630
P	167	HIS	-	EXPRESSION TAG	UNP Q16630
P	168	HIS	-	EXPRESSION TAG	UNP Q16630
P	169	HIS	-	EXPRESSION TAG	UNP Q16630
Q	159	SER	CYS	ENGINEERED MUTATION	UNP Q16630
Q	162	LEU	-	EXPRESSION TAG	UNP Q16630
Q	163	GLU	-	EXPRESSION TAG	UNP Q16630
Q	164	HIS	-	EXPRESSION TAG	UNP Q16630
Q	165	HIS	-	EXPRESSION TAG	UNP Q16630
Q	166	HIS	-	EXPRESSION TAG	UNP Q16630
Q	167	HIS	-	EXPRESSION TAG	UNP Q16630
Q	168	HIS	-	EXPRESSION TAG	UNP Q16630
Q	169	HIS	-	EXPRESSION TAG	UNP Q16630

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	30	Total O 30 30	0	0
3	B	40	Total O 40 40	0	0
3	C	23	Total O 23 23	0	0
3	D	49	Total O 49 49	0	0
3	E	13	Total O 13 13	0	0
3	F	27	Total O 27 27	0	0
3	L	3	Total O 3 3	0	0
3	M	16	Total O 16 16	0	0
3	N	21	Total O 21 21	0	0

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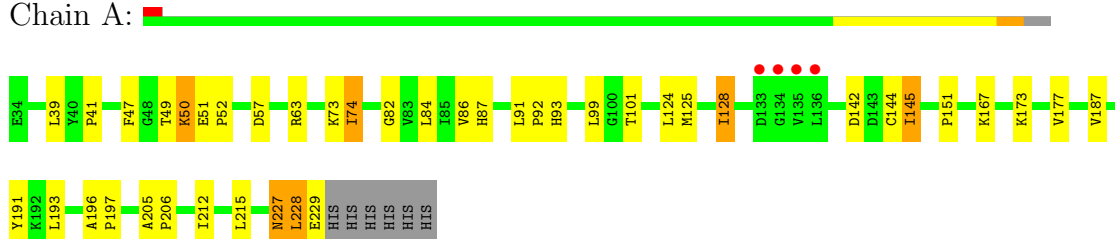
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	O	12	Total 12	O 12	0	0
3	P	6	Total 6	O 6	0	0
3	Q	8	Total 8	O 8	0	0

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.

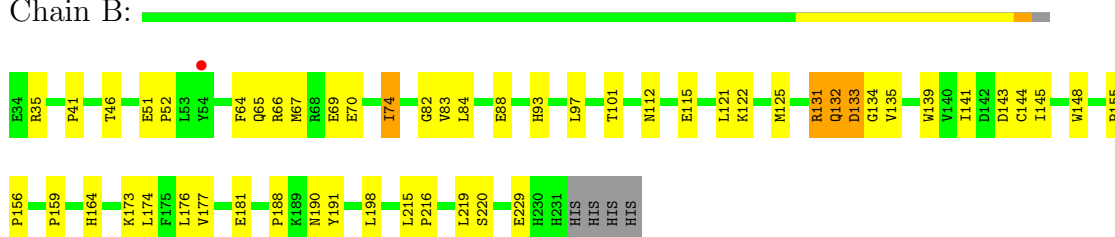
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain A:



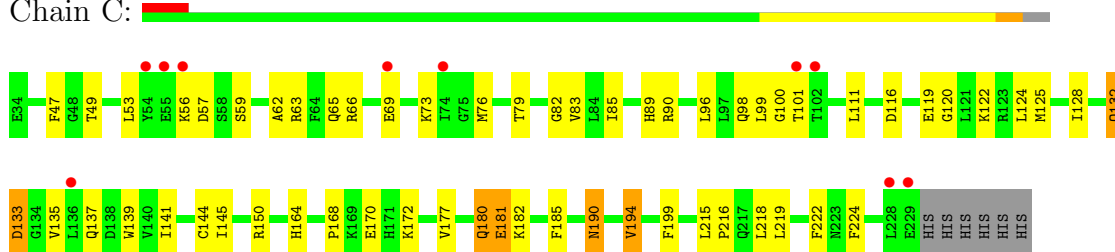
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain B:



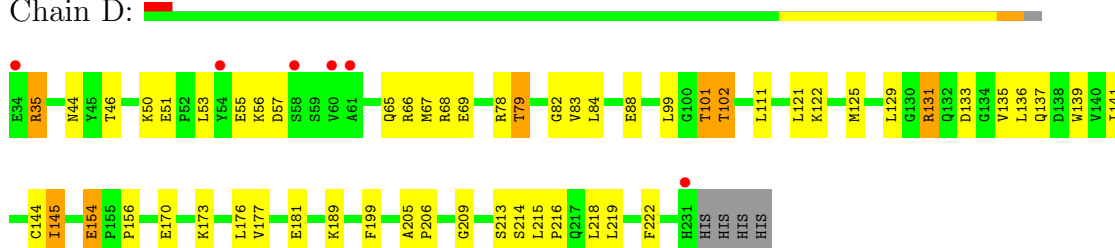
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain C:



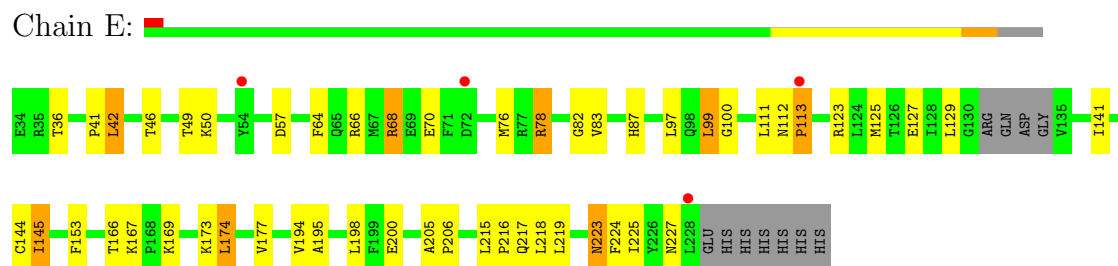
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain D:



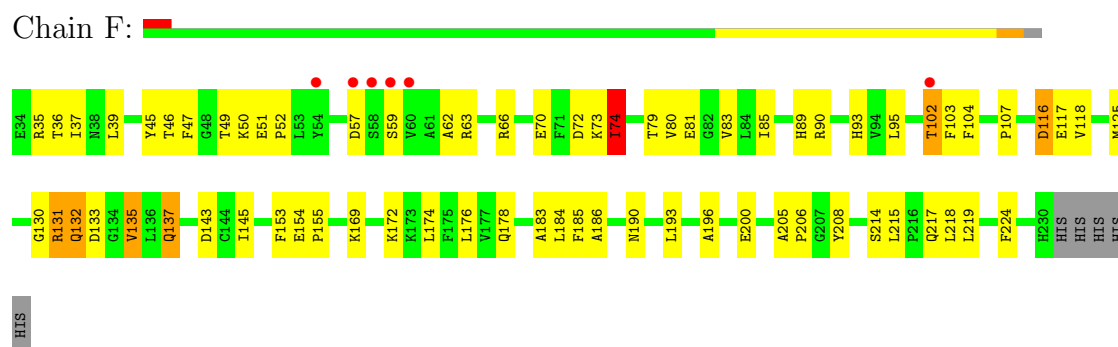
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain E:



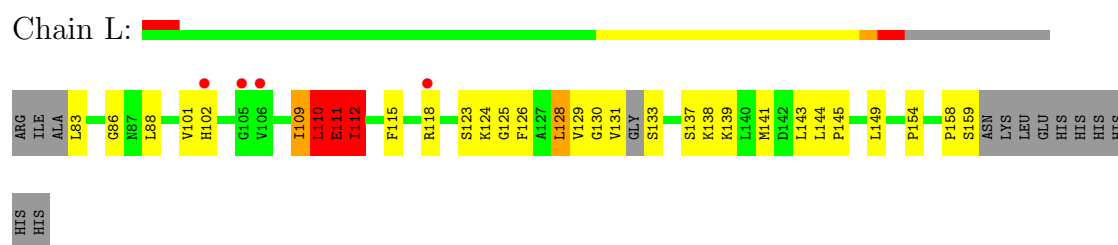
- Molecule 1: Cleavage and polyadenylation specificity factor subunit 5

Chain F:



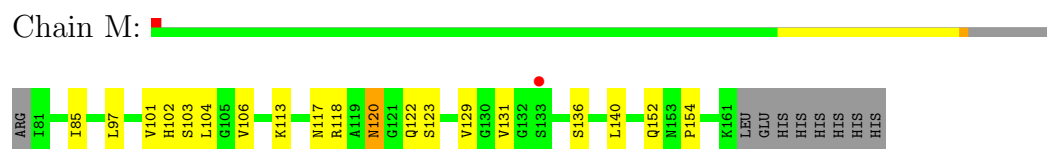
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain L:



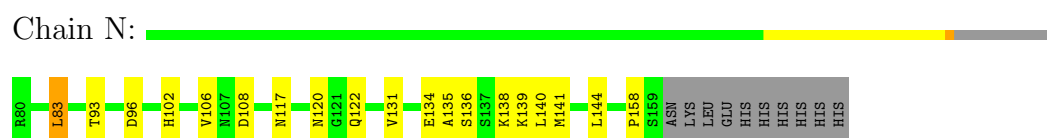
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain M:



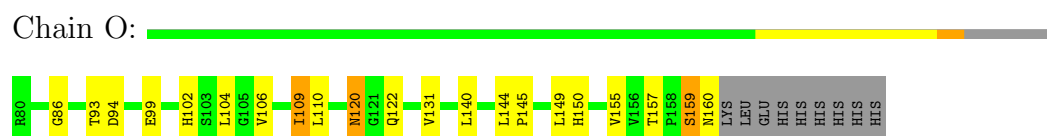
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain N:



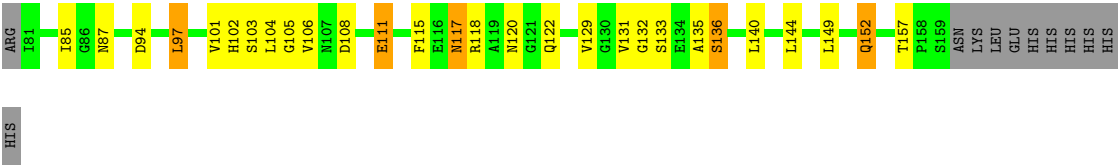
- Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain O:



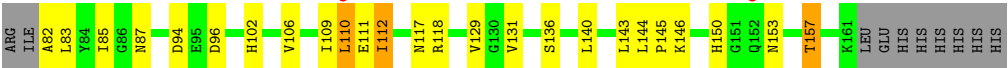
● Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain P: 



● Molecule 2: Cleavage and polyadenylation specificity factor subunit 6

Chain Q: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	160.44Å 105.69Å 147.08Å 90.00° 112.72° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 29.94 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.9 (15.00-2.70) 98.9 (29.94-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.29 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.214 , 0.265 0.214 , 0.267	Depositor DCC
R_{free} test set	3119 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 16.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 61654 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13253	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.49	0/1635	0.62	1/2223 (0.0%)
1	B	0.50	0/1658	0.63	0/2255
1	C	0.44	0/1620	0.62	0/2205
1	D	0.53	0/1661	0.67	1/2259 (0.0%)
1	E	0.43	0/1564	0.58	0/2133
1	F	0.45	0/1644	0.61	0/2238
2	L	0.57	0/537	0.97	4/735 (0.5%)
2	M	0.52	0/615	0.63	0/839
2	N	0.53	0/616	0.68	0/838
2	O	0.56	0/611	0.72	2/835 (0.2%)
2	P	0.51	0/598	0.65	0/816
2	Q	0.56	0/595	0.67	0/812
All	All	0.49	0/13354	0.65	8/18188 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	F	0	1
2	L	1	2
All	All	1	3

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	112	ILE	CG1-CB-CG2	9.27	131.78	111.40
2	L	110	LEU	O-C-N	-9.20	107.98	122.70
2	L	112	ILE	CB-CA-C	-8.97	93.66	111.60
2	O	109	ILE	CB-CA-C	-7.03	97.55	111.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	219	LEU	CA-CB-CG	5.76	128.56	115.30
2	O	120	ASN	CB-CA-C	-5.60	99.20	110.40
1	A	74	ILE	CB-CA-C	-5.10	101.40	111.60
2	L	111	GLU	O-C-N	-5.01	114.68	122.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	L	112	ILE	CB

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	132	GLN	Peptide
2	L	110	LEU	Mainchain
2	L	111	GLU	Mainchain

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	1589	0	1578	33	0
1	B	1611	0	1586	56	0
1	C	1574	0	1547	54	0
1	D	1613	0	1594	50	0
1	E	1519	0	1471	47	0
1	F	1597	0	1563	69	0
2	L	527	0	446	32	0
2	M	603	0	573	11	0
2	N	604	0	589	14	0
2	O	599	0	564	13	0
2	P	586	0	553	33	0
2	Q	583	0	536	27	0
3	A	30	0	0	1	0
3	B	40	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	23	0	0	0	0
3	D	49	0	0	1	0
3	E	13	0	0	1	0
3	F	27	0	0	0	0
3	L	3	0	0	0	0
3	M	16	0	0	0	0
3	N	21	0	0	0	0
3	O	12	0	0	0	0
3	P	6	0	0	0	0
3	Q	8	0	0	0	0
All	All	13253	0	12600	427	0

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 17.

All (427) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:128:LEU:HD12	2:L:129:VAL:N	1.27	1.41
2:P:117:ASN:ND2	2:P:120:ASN:HB2	1.40	1.36
2:L:128:LEU:HD12	2:L:128:LEU:C	1.47	1.21
2:L:110:LEU:N	2:L:130:GLY:O	1.71	1.21
2:P:117:ASN:HD21	2:P:120:ASN:CB	1.55	1.17
1:D:35:ARG:HH11	1:D:35:ARG:CG	1.59	1.16
1:E:99:LEU:C	1:E:99:LEU:HD23	1.65	1.14
1:B:132:GLN:HA	1:B:132:GLN:OE1	1.44	1.13
1:D:35:ARG:NH1	1:D:35:ARG:HG2	1.50	1.12
2:L:112:ILE:HG23	2:L:129:VAL:HG22	1.33	1.09
1:C:132:GLN:OE1	1:C:132:GLN:HA	1.49	1.07
2:L:128:LEU:CD1	2:L:129:VAL:N	2.20	1.05
1:F:74:ILE:O	1:F:74:ILE:HG12	1.58	1.03
1:F:131:ARG:NH1	1:F:135:VAL:HB	1.74	1.02
2:L:112:ILE:CG2	2:L:129:VAL:HG22	1.93	0.98
1:A:228:LEU:O	1:A:229:GLU:CB	2.09	0.97
1:F:131:ARG:NH1	1:F:135:VAL:CG2	2.28	0.96
1:C:83:VAL:HG21	1:C:145:ILE:HD11	1.48	0.96
1:B:74:ILE:HG13	1:B:74:ILE:O	1.65	0.95
1:F:131:ARG:NH1	1:F:135:VAL:CB	2.31	0.94
2:Q:110:LEU:CD2	2:Q:110:LEU:C	2.34	0.94
1:D:65:GLN:NE2	1:D:69:GLU:OE2	2.01	0.94
1:F:131:ARG:HH12	1:F:135:VAL:HG21	1.32	0.91
1:B:65:GLN:O	1:B:69:GLU:HG3	1.71	0.91
1:B:66:ARG:HD2	1:B:70:GLU:OE2	1.71	0.90

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:99:LEU:HD23	1:E:100:GLY:N	1.86	0.90
1:D:35:ARG:HH11	1:D:35:ARG:HG2	0.75	0.90
1:E:99:LEU:C	1:E:99:LEU:CD2	2.38	0.89
1:E:125:MET:HE3	1:E:177:VAL:HG21	1.52	0.89
2:L:138:LYS:HA	2:L:141:MET:CE	2.02	0.89
1:C:57:ASP:OD1	1:C:63:ARG:HA	1.73	0.88
1:F:66:ARG:O	1:F:70:GLU:HG3	1.72	0.87
2:L:128:LEU:CD1	2:L:128:LEU:C	2.29	0.87
1:F:47:PHE:CE1	1:F:185:PHE:HD2	1.92	0.86
2:Q:82:ALA:O	2:Q:83:LEU:HD12	1.75	0.86
1:A:99:LEU:HD11	1:A:191:TYR:HE2	1.41	0.85
1:F:131:ARG:NH1	1:F:135:VAL:HG21	1.89	0.85
1:A:99:LEU:HD11	1:A:191:TYR:CE2	2.12	0.85
1:B:132:GLN:O	1:B:134:GLY:N	2.09	0.84
1:E:144:CYS:SG	1:E:173:LYS:HE3	2.18	0.83
1:F:131:ARG:HH11	1:F:135:VAL:HB	1.43	0.83
1:F:46:THR:OG1	1:F:184:LEU:HD12	1.80	0.82
1:C:56:LYS:HD3	1:C:56:LYS:N	1.95	0.81
1:A:227:ASN:H	1:A:227:ASN:HD22	1.29	0.80
1:F:74:ILE:O	1:F:74:ILE:CG1	2.29	0.79
2:Q:110:LEU:C	2:Q:110:LEU:HD23	2.01	0.79
1:A:227:ASN:ND2	1:A:227:ASN:H	1.79	0.78
1:B:132:GLN:OE1	1:B:132:GLN:CA	2.30	0.78
2:L:128:LEU:HD12	2:L:129:VAL:H	1.46	0.77
2:Q:110:LEU:C	2:Q:110:LEU:HD22	2.02	0.77
1:B:74:ILE:CG1	1:B:74:ILE:O	2.32	0.77
2:L:138:LYS:HA	2:L:141:MET:HE2	1.67	0.76
2:Q:118:ARG:O	2:Q:118:ARG:HG2	1.84	0.75
1:F:39:LEU:HD11	1:F:224:PHE:CD1	2.22	0.75
1:F:49:THR:CG2	1:F:49:THR:O	2.36	0.74
1:A:144:CYS:SG	1:A:173:LYS:HE3	2.28	0.74
1:C:132:GLN:OE1	1:C:132:GLN:CA	2.29	0.73
1:F:47:PHE:CE1	1:F:185:PHE:CD2	2.76	0.73
2:N:135:ALA:O	2:N:139:LYS:HG3	1.88	0.73
1:C:124:LEU:O	1:C:128:ILE:HG12	1.89	0.72
1:B:132:GLN:C	1:B:134:GLY:N	2.42	0.72
1:B:83:VAL:HG21	1:B:145:ILE:CD1	2.20	0.71
1:F:46:THR:OG1	1:F:184:LEU:CD1	2.38	0.71
2:Q:110:LEU:O	2:Q:110:LEU:HD23	1.90	0.71
1:C:56:LYS:HD3	1:C:56:LYS:H	1.56	0.70
1:F:131:ARG:HH11	1:F:135:VAL:CG2	2.03	0.70
2:M:120:ASN:HB3	2:M:122:GLN:H	1.57	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:133:ASP:C	1:C:133:ASP:OD1	2.28	0.70
2:Q:82:ALA:C	2:Q:83:LEU:HD12	2.11	0.70
1:B:141:ILE:HG12	1:B:177:VAL:HG22	1.72	0.69
2:L:144:LEU:HB3	2:L:145:PRO:HD3	1.74	0.69
1:B:132:GLN:C	1:B:134:GLY:H	1.94	0.69
1:E:205:ALA:HB3	1:E:206:PRO:HD3	1.74	0.69
2:L:110:LEU:CB	2:L:130:GLY:O	2.41	0.69
2:L:138:LYS:HA	2:L:141:MET:HE1	1.73	0.69
1:B:131:ARG:NH1	1:B:133:ASP:OD2	2.27	0.68
1:D:144:CYS:SG	1:D:173:LYS:HE3	2.33	0.68
1:F:131:ARG:HH11	1:F:135:VAL:CB	2.02	0.68
2:P:133:SER:OG	2:P:136:SER:OG	2.12	0.67
1:C:124:LEU:O	1:C:128:ILE:CG1	2.42	0.67
1:D:136:LEU:HG	1:D:137:GLN:N	2.09	0.67
2:Q:106:VAL:HG13	2:Q:136:SER:HB3	1.76	0.67
2:P:87:ASN:O	2:P:152:GLN:NE2	2.27	0.67
2:Q:143:LEU:O	2:Q:146:LYS:CB	2.43	0.67
2:L:110:LEU:CA	2:L:130:GLY:O	2.42	0.67
2:P:111:GLU:N	2:P:111:GLU:OE2	2.28	0.66
1:F:83:VAL:HG21	1:F:145:ILE:HD11	1.76	0.66
1:C:65:GLN:O	1:C:69:GLU:HG3	1.95	0.66
1:B:66:ARG:CD	1:B:70:GLU:OE2	2.42	0.66
2:P:117:ASN:ND2	2:P:117:ASN:O	2.27	0.66
1:F:116:ASP:OD2	1:F:117:GLU:N	2.29	0.66
1:F:131:ARG:NH1	1:F:133:ASP:OD2	2.28	0.66
1:E:82:GLY:HA3	1:E:125:MET:CE	2.26	0.65
1:E:78:ARG:HB3	1:E:111:LEU:HD12	1.78	0.65
1:D:56:LYS:HG2	1:D:56:LYS:O	1.97	0.65
1:F:116:ASP:OD2	1:F:116:ASP:C	2.34	0.65
2:L:101:VAL:HG11	2:L:109:ILE:HD13	1.79	0.65
1:F:215:LEU:O	1:F:219:LEU:HG	1.96	0.65
2:Q:96:ASP:OD2	2:Q:150:HIS:NE2	2.26	0.65
1:F:153:PHE:O	1:F:154:GLU:HG2	1.96	0.64
1:F:57:ASP:OD2	1:F:66:ARG:NH1	2.22	0.64
1:B:83:VAL:HG21	1:B:145:ILE:HD11	1.78	0.64
2:O:109:ILE:O	2:O:109:ILE:HG22	1.93	0.64
1:C:98:GLN:HG2	1:C:99:LEU:N	2.13	0.64
1:C:125:MET:HE3	1:C:177:VAL:HG21	1.79	0.64
1:B:131:ARG:NH1	1:B:135:VAL:CG2	2.61	0.64
1:E:42:LEU:O	1:E:42:LEU:HD12	1.97	0.63
1:F:153:PHE:CZ	1:F:169:LYS:HD2	2.33	0.63
2:L:137:SER:O	2:L:141:MET:HG3	1.99	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:217:GLN:OE1	1:E:217:GLN:N	2.31	0.63
2:N:120:ASN:HB3	2:N:122:GLN:H	1.64	0.63
2:N:141:MET:SD	2:N:158:PRO:HD3	2.39	0.63
2:P:117:ASN:ND2	2:P:120:ASN:CB	2.31	0.62
2:Q:85:ILE:HD12	2:Q:129:VAL:HG21	1.80	0.62
1:A:74:ILE:HG22	1:A:74:ILE:O	2.00	0.62
2:L:110:LEU:CB	2:L:130:GLY:C	2.68	0.62
1:D:131:ARG:HD3	1:D:133:ASP:OD1	1.99	0.62
1:E:66:ARG:NH2	1:E:70:GLU:OE2	2.32	0.62
1:E:82:GLY:HA3	1:E:125:MET:HE1	1.82	0.61
1:C:133:ASP:OD1	1:C:135:VAL:N	2.23	0.61
1:B:82:GLY:HA3	1:B:125:MET:HE1	1.81	0.61
2:Q:140:LEU:HD22	2:Q:144:LEU:HD22	1.82	0.61
1:C:59:SER:H	1:C:62:ALA:HB3	1.66	0.61
2:P:140:LEU:HD22	2:P:144:LEU:HD22	1.82	0.60
1:D:199:PHE:HB2	2:O:120:ASN:HB3	1.83	0.60
1:F:214:SER:HA	1:F:217:GLN:OE1	2.01	0.60
1:B:144:CYS:SG	1:B:173:LYS:HE3	2.40	0.60
1:B:215:LEU:O	1:B:219:LEU:HG	2.01	0.60
1:C:76:MET:HG3	1:C:168:PRO:O	2.02	0.60
1:F:131:ARG:HH12	1:F:135:VAL:CG2	1.97	0.60
1:C:49:THR:O	1:C:49:THR:CG2	2.49	0.60
1:C:49:THR:O	1:C:49:THR:HG23	2.01	0.60
2:M:85:ILE:HD12	2:M:129:VAL:HG21	1.84	0.60
1:D:122:LYS:HG3	1:D:141:ILE:HD12	1.84	0.60
1:D:50:LYS:HG3	1:D:51:GLU:N	2.16	0.60
1:F:45:TYR:HD1	1:F:183:ALA:HB3	1.67	0.59
2:O:120:ASN:HB2	2:O:122:GLN:H	1.68	0.59
2:O:94:ASP:N	2:O:94:ASP:OD1	2.34	0.59
1:C:133:ASP:OD1	1:C:135:VAL:HB	2.03	0.58
1:C:190:ASN:OD1	1:C:190:ASN:N	2.32	0.58
1:F:102:THR:HG23	1:F:102:THR:O	2.02	0.58
2:N:83:LEU:HD12	2:N:158:PRO:HA	1.86	0.58
1:A:227:ASN:ND2	1:A:227:ASN:N	2.50	0.58
1:B:135:VAL:HG13	2:P:115:PHE:HZ	1.69	0.57
1:E:83:VAL:HG21	1:E:145:ILE:HD12	1.86	0.57
1:B:83:VAL:HG21	1:B:145:ILE:HD13	1.87	0.57
2:P:117:ASN:HD21	2:P:120:ASN:HB2	0.61	0.57
1:B:66:ARG:CG	1:B:70:GLU:OE2	2.53	0.57
1:B:65:GLN:O	1:B:69:GLU:CG	2.50	0.57
1:E:125:MET:CE	1:E:177:VAL:HG21	2.28	0.57
1:A:151:PRO:HD2	3:A:244:HOH:O	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:74:ILE:O	1:A:74:ILE:CG2	2.53	0.57
1:E:83:VAL:HG21	1:E:145:ILE:CD1	2.35	0.57
1:F:89:HIS:O	1:F:90:ARG:HB2	2.05	0.56
1:C:56:LYS:CD	1:C:56:LYS:N	2.57	0.56
1:E:123:ARG:NH1	1:E:127:GLU:OE2	2.38	0.56
2:Q:112:ILE:HG12	2:Q:112:ILE:O	2.05	0.56
2:P:102:HIS:O	2:P:105:GLY:N	2.25	0.56
2:L:149:LEU:HD12	2:L:154:PRO:HG3	1.87	0.56
1:D:205:ALA:N	1:D:206:PRO:HD2	2.20	0.56
2:Q:118:ARG:O	2:Q:118:ARG:CG	2.51	0.56
2:P:111:GLU:O	2:P:111:GLU:OE2	2.23	0.56
1:E:41:PRO:HD3	1:E:227:ASN:O	2.05	0.56
1:B:135:VAL:HG13	2:P:115:PHE:CZ	2.41	0.56
1:C:101:THR:HG22	1:C:101:THR:O	2.05	0.56
2:P:106:VAL:HG13	2:P:136:SER:HB3	1.86	0.55
2:Q:102:HIS:HD2	2:Q:106:VAL:O	1.89	0.55
2:Q:94:ASP:OD1	2:Q:94:ASP:N	2.39	0.55
1:D:136:LEU:HG	1:D:137:GLN:H	1.71	0.55
1:F:131:ARG:HD3	1:F:133:ASP:CB	2.37	0.55
2:Q:110:LEU:CD2	2:Q:110:LEU:O	2.52	0.55
1:F:95:LEU:HD13	1:F:193:LEU:HD21	1.89	0.54
1:B:148:TRP:CZ3	1:B:159:PRO:HD3	2.42	0.54
1:C:85:ILE:HD11	1:C:219:LEU:HD22	1.89	0.54
1:D:101:THR:HG22	1:D:102:THR:HG22	1.89	0.54
1:D:55:GLU:C	1:D:57:ASP:H	2.09	0.54
1:D:214:SER:O	1:D:218:LEU:HG	2.07	0.54
2:L:112:ILE:HG23	2:L:129:VAL:CG2	2.22	0.54
1:E:145:ILE:HG13	1:E:174:LEU:HB3	1.89	0.54
2:M:106:VAL:HG13	2:M:136:SER:HB3	1.89	0.54
1:D:205:ALA:N	1:D:206:PRO:CD	2.70	0.54
1:F:49:THR:O	1:F:49:THR:HG23	2.03	0.54
1:D:83:VAL:HG21	1:D:145:ILE:HD11	1.89	0.54
2:Q:144:LEU:HB3	2:Q:145:PRO:HD3	1.89	0.54
2:P:111:GLU:OE2	2:P:111:GLU:CA	2.56	0.53
1:D:79:THR:OG1	1:D:170:GLU:OE2	2.21	0.53
1:D:56:LYS:O	1:D:56:LYS:CG	2.56	0.53
1:B:155:PRO:HB2	1:B:156:PRO:HD3	1.90	0.53
1:F:49:THR:O	1:F:49:THR:HG22	2.09	0.53
1:D:215:LEU:N	1:D:216:PRO:CD	2.71	0.53
1:F:205:ALA:N	1:F:206:PRO:CD	2.72	0.53
1:B:131:ARG:NH1	1:B:135:VAL:HG23	2.23	0.52
1:B:82:GLY:HA3	1:B:125:MET:CE	2.40	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:98:GLN:NE2	1:C:100:GLY:O	2.41	0.52
1:C:164:HIS:HA	2:O:93:THR:HG22	1.92	0.52
1:C:116:ASP:HB3	1:C:119:GLU:HB2	1.92	0.52
1:B:198:LEU:HD21	1:B:219:LEU:HD12	1.92	0.52
1:C:66:ARG:O	1:C:66:ARG:CG	2.58	0.52
2:Q:82:ALA:C	2:Q:83:LEU:CD1	2.77	0.51
1:A:49:THR:HA	1:A:187:VAL:O	2.10	0.51
1:B:133:ASP:C	1:B:133:ASP:OD1	2.47	0.51
2:L:158:PRO:O	2:L:159:SER:C	2.48	0.51
1:D:209:GLY:O	1:D:213:SER:OG	2.19	0.51
2:Q:144:LEU:C	2:Q:146:LYS:H	2.13	0.51
1:D:101:THR:CG2	1:D:101:THR:O	2.54	0.51
2:P:97:LEU:O	2:P:101:VAL:HG23	2.11	0.51
1:B:131:ARG:NH1	1:B:133:ASP:CG	2.64	0.51
1:E:99:LEU:O	1:E:99:LEU:CD2	2.58	0.51
1:C:122:LYS:HG3	1:C:141:ILE:HD12	1.93	0.51
1:E:64:PHE:O	1:E:68:ARG:HG2	2.11	0.51
2:M:101:VAL:HG13	2:M:140:LEU:HD21	1.93	0.51
1:F:52:PRO:HA	1:F:190:ASN:OD1	2.11	0.51
1:A:124:LEU:O	1:A:128:ILE:HG12	2.11	0.51
2:P:102:HIS:HD2	2:P:106:VAL:O	1.93	0.51
1:E:200:GLU:O	1:E:200:GLU:HG3	2.11	0.51
1:D:141:ILE:HG12	1:D:177:VAL:HG22	1.94	0.50
1:E:83:VAL:CG2	1:E:145:ILE:HD11	2.42	0.50
2:O:159:SER:O	2:O:160:ASN:OD1	2.29	0.50
1:B:35:ARG:NH1	1:B:35:ARG:HG3	2.26	0.50
1:E:49:THR:HG22	1:E:50:LYS:N	2.26	0.50
2:P:117:ASN:C	2:P:117:ASN:ND2	2.64	0.50
1:C:124:LEU:O	1:C:128:ILE:HG13	2.12	0.50
2:N:102:HIS:HD2	2:N:106:VAL:O	1.95	0.50
2:Q:110:LEU:HD22	2:Q:111:GLU:N	2.26	0.50
1:F:81:GLU:HB2	1:F:174:LEU:HD22	1.94	0.50
1:A:124:LEU:O	1:A:128:ILE:CG1	2.60	0.50
1:F:103:PHE:C	1:F:103:PHE:CD1	2.83	0.50
2:P:111:GLU:N	2:P:111:GLU:CD	2.65	0.50
1:A:50:LYS:HE3	1:A:128:ILE:O	2.12	0.50
1:E:99:LEU:O	1:E:99:LEU:HD23	2.10	0.50
1:B:132:GLN:O	1:B:133:ASP:C	2.50	0.50
1:F:130:GLY:HA2	1:F:137:GLN:OE1	2.11	0.49
2:P:101:VAL:HG12	2:P:106:VAL:HB	1.94	0.49
1:F:116:ASP:OD2	1:F:118:VAL:N	2.45	0.49
1:E:66:ARG:HG2	1:E:66:ARG:NH2	2.28	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:68:ARG:O	1:D:68:ARG:HG3	2.12	0.49
2:P:94:ASP:N	2:P:94:ASP:OD1	2.43	0.49
2:N:108:ASP:N	2:N:108:ASP:OD1	2.40	0.49
2:L:86:GLY:HA2	2:L:126:PHE:HB3	1.93	0.49
2:L:144:LEU:N	2:L:145:PRO:CD	2.76	0.49
2:M:102:HIS:C	2:M:104:LEU:H	2.16	0.49
1:C:215:LEU:N	1:C:216:PRO:CD	2.76	0.49
1:F:47:PHE:HE1	1:F:185:PHE:CD2	2.30	0.49
2:L:102:HIS:CD2	2:L:109:ILE:HD11	2.47	0.49
1:F:83:VAL:HG21	1:F:145:ILE:CD1	2.42	0.48
1:D:122:LYS:HG2	1:D:139:TRP:HB2	1.93	0.48
1:F:103:PHE:HD1	1:F:104:PHE:N	2.10	0.48
1:E:46:THR:HG23	3:E:7:HOH:O	2.13	0.48
1:D:215:LEU:HD23	1:D:218:LEU:HD12	1.95	0.48
1:F:103:PHE:CD1	1:F:104:PHE:N	2.81	0.48
1:E:82:GLY:HA3	1:E:125:MET:HE2	1.96	0.48
1:A:205:ALA:N	1:A:206:PRO:CD	2.76	0.48
1:B:131:ARG:HH11	1:B:133:ASP:CG	2.15	0.48
1:E:66:ARG:HH21	1:E:66:ARG:CG	2.26	0.48
1:B:35:ARG:HH11	1:B:35:ARG:HG3	1.78	0.48
1:D:88:GLU:HG2	3:D:258:HOH:O	2.13	0.48
1:E:215:LEU:O	1:E:219:LEU:HG	2.13	0.48
2:P:149:LEU:HA	2:P:149:LEU:HD23	1.57	0.48
2:O:102:HIS:HA	2:O:106:VAL:O	2.13	0.48
2:P:111:GLU:C	2:P:111:GLU:OE2	2.52	0.48
1:E:205:ALA:HB3	1:E:206:PRO:CD	2.44	0.48
2:L:86:GLY:HA2	2:L:126:PHE:CB	2.44	0.48
1:F:200:GLU:OE2	1:F:208:TYR:OH	2.21	0.48
2:L:111:GLU:O	2:L:130:GLY:N	2.47	0.48
1:D:78:ARG:HB3	1:D:111:LEU:HD12	1.95	0.48
1:A:82:GLY:HA3	1:A:125:MET:CE	2.44	0.48
2:N:134:GLU:OE2	2:N:138:LYS:HE3	2.13	0.47
2:N:131:VAL:CG1	2:N:136:SER:HB2	2.44	0.47
1:F:131:ARG:HH12	1:F:135:VAL:CB	2.22	0.47
1:C:122:LYS:HG2	1:C:139:TRP:HB2	1.96	0.47
2:N:131:VAL:HG13	2:N:136:SER:HB2	1.97	0.47
1:D:122:LYS:HE2	1:D:139:TRP:O	2.15	0.47
1:D:101:THR:HG23	1:D:101:THR:O	2.15	0.47
1:D:66:ARG:C	1:D:68:ARG:N	2.64	0.47
2:O:86:GLY:N	2:O:155:VAL:O	2.42	0.47
1:D:82:GLY:HA3	1:D:125:MET:CE	2.44	0.47
1:B:131:ARG:NH1	1:B:133:ASP:OD1	2.42	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:56:LYS:CD	1:C:56:LYS:H	2.19	0.47
1:F:93:HIS:HA	1:F:196:ALA:O	2.15	0.47
2:M:152:GLN:O	2:M:154:PRO:HD3	2.15	0.47
1:B:121:LEU:O	1:B:125:MET:HG2	2.15	0.46
2:M:101:VAL:HG12	2:M:106:VAL:HB	1.96	0.46
2:O:144:LEU:HB3	2:O:145:PRO:HD3	1.97	0.46
1:A:47:PHE:CZ	1:A:193:LEU:HD23	2.50	0.46
2:N:117:ASN:OD1	2:N:120:ASN:HB2	2.16	0.46
1:F:45:TYR:CD1	1:F:183:ALA:HB3	2.49	0.46
1:F:79:THR:CG2	1:F:172:LYS:HG3	2.45	0.46
2:N:140:LEU:HD22	2:N:144:LEU:HD22	1.97	0.46
1:E:153:PHE:CZ	1:E:169:LYS:HE2	2.50	0.46
1:A:73:LYS:NZ	2:Q:157:THR:HB	2.29	0.46
1:B:88:GLU:HB2	1:B:93:HIS:CE1	2.51	0.46
1:C:98:GLN:CG	1:C:99:LEU:N	2.78	0.46
1:E:112:ASN:O	1:E:113:PRO:C	2.54	0.46
1:E:224:PHE:O	1:E:225:ILE:HD13	2.16	0.46
2:O:99:GLU:O	2:O:99:GLU:HG2	2.15	0.46
1:A:41:PRO:HA	1:A:87:HIS:O	2.15	0.46
1:B:66:ARG:O	1:B:70:GLU:HG3	2.15	0.46
1:D:66:ARG:C	1:D:68:ARG:H	2.19	0.46
1:C:215:LEU:N	1:C:216:PRO:HD2	2.31	0.46
1:B:181:GLU:HB3	2:P:118:ARG:HD2	1.98	0.46
2:Q:87:ASN:ND2	2:Q:153:ASN:O	2.44	0.46
1:E:205:ALA:N	1:E:206:PRO:HD2	2.31	0.45
1:C:47:PHE:CE1	1:C:185:PHE:HD2	2.34	0.45
1:F:154:GLU:HB3	1:F:155:PRO:HD2	1.99	0.45
1:E:153:PHE:CE2	1:E:169:LYS:HE2	2.52	0.45
2:L:139:LYS:O	2:L:143:LEU:CB	2.64	0.45
1:D:129:LEU:HD23	1:D:129:LEU:HA	1.75	0.45
1:C:125:MET:CE	1:C:177:VAL:HG21	2.46	0.45
1:F:102:THR:CG2	1:F:102:THR:O	2.63	0.45
1:C:85:ILE:HD11	1:C:219:LEU:CD2	2.46	0.45
1:D:78:ARG:O	1:D:111:LEU:HB2	2.17	0.45
2:P:120:ASN:HB3	2:P:122:GLN:H	1.82	0.45
1:D:176:LEU:HD12	1:D:222:PHE:CZ	2.52	0.45
1:B:122:LYS:HG2	1:B:139:TRP:HB2	1.99	0.44
2:P:102:HIS:C	2:P:104:LEU:H	2.20	0.44
1:D:66:ARG:O	1:D:68:ARG:N	2.50	0.44
1:D:44:ASN:HB3	1:D:181:GLU:O	2.17	0.44
2:P:102:HIS:C	2:P:104:LEU:N	2.71	0.44
1:A:47:PHE:CD1	1:A:47:PHE:N	2.85	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:176:LEU:HD23	1:F:176:LEU:C	2.38	0.44
2:P:85:ILE:HD12	2:P:129:VAL:HG21	1.98	0.44
1:F:104:PHE:N	1:F:104:PHE:CD2	2.86	0.44
2:Q:85:ILE:HD12	2:Q:129:VAL:CG2	2.45	0.44
2:O:149:LEU:O	2:O:150:HIS:C	2.56	0.44
1:B:125:MET:HB3	1:B:139:TRP:CZ2	2.53	0.44
2:L:115:PHE:O	2:L:124:LYS:HB2	2.17	0.44
1:E:198:LEU:HD21	1:E:219:LEU:HD12	2.00	0.44
2:L:83:LEU:N	2:L:129:VAL:O	2.50	0.43
1:A:39:LEU:O	1:A:227:ASN:ND2	2.46	0.43
1:C:69:GLU:O	1:C:73:LYS:HG3	2.18	0.43
1:A:99:LEU:HD11	1:A:191:TYR:CD2	2.51	0.43
1:A:51:GLU:HB3	1:A:52:PRO:HD2	2.00	0.43
1:B:131:ARG:CZ	1:B:135:VAL:HB	2.48	0.43
2:M:117:ASN:HB2	2:M:120:ASN:HB2	2.01	0.43
1:D:154:GLU:HB3	1:D:156:PRO:HD2	2.01	0.43
1:E:218:LEU:HD23	1:F:218:LEU:HD23	2.00	0.43
2:M:97:LEU:HD23	2:M:97:LEU:HA	1.82	0.43
2:P:108:ASP:OD2	2:P:132:GLY:CA	2.66	0.43
1:B:215:LEU:N	1:B:216:PRO:CD	2.82	0.43
1:F:51:GLU:OE1	1:F:52:PRO:HD2	2.19	0.43
1:B:64:PHE:HD1	1:B:67:MET:CE	2.32	0.43
2:L:128:LEU:CD1	2:L:129:VAL:H	2.14	0.43
1:C:111:LEU:CD2	1:C:120:GLY:HA3	2.49	0.43
1:E:68:ARG:HA	1:E:153:PHE:CE1	2.54	0.43
1:A:212:ILE:HG23	1:A:215:LEU:HD12	1.99	0.43
1:B:188:PRO:O	1:B:190:ASN:N	2.51	0.43
1:C:180:GLN:O	1:C:181:GLU:C	2.57	0.43
1:E:76:MET:HB3	1:E:76:MET:HE3	1.83	0.43
1:B:66:ARG:HG2	1:B:70:GLU:CD	2.39	0.43
1:F:145:ILE:HG13	1:F:174:LEU:HB2	2.00	0.43
1:B:51:GLU:HB3	1:B:52:PRO:HD2	1.99	0.43
1:B:164:HIS:CG	2:L:123:SER:HB2	2.54	0.43
1:F:131:ARG:HD3	1:F:133:ASP:HB2	2.00	0.42
2:L:102:HIS:NE2	2:L:109:ILE:HD11	2.33	0.42
2:L:88:LEU:HB2	2:L:125:GLY:O	2.19	0.42
1:E:166:THR:OG1	1:E:167:LYS:HE3	2.19	0.42
1:F:59:SER:O	1:F:62:ALA:HB3	2.19	0.42
1:E:66:ARG:CG	1:E:66:ARG:NH2	2.81	0.42
1:D:55:GLU:C	1:D:57:ASP:N	2.73	0.42
2:Q:87:ASN:CG	2:Q:87:ASN:O	2.57	0.42
1:C:96:LEU:HB2	1:C:194:VAL:HG23	2.00	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:107:PRO:HB2	1:F:125:MET:SD	2.59	0.42
1:B:41:PRO:HG2	1:B:229:GLU:HG2	2.00	0.42
1:B:112:ASN:O	1:B:115:GLU:HB3	2.19	0.42
1:C:98:GLN:HG2	1:C:99:LEU:H	1.82	0.42
1:C:82:GLY:HA3	1:C:125:MET:CE	2.49	0.42
1:E:83:VAL:CG2	1:E:145:ILE:CD1	2.98	0.42
1:D:82:GLY:HA3	1:D:125:MET:HE1	2.02	0.42
1:A:47:PHE:HD1	1:A:47:PHE:N	2.17	0.42
1:B:143:ASP:HB2	1:B:176:LEU:HB3	2.01	0.42
1:C:150:ARG:NE	1:C:170:GLU:OE1	2.44	0.42
1:A:84:LEU:HD12	1:A:177:VAL:HB	2.02	0.42
2:P:133:SER:C	2:P:135:ALA:N	2.73	0.42
1:F:36:THR:C	1:F:37:ILE:HG12	2.39	0.42
2:P:133:SER:C	2:P:135:ALA:H	2.22	0.42
1:C:141:ILE:HG12	1:C:177:VAL:HG22	2.02	0.42
1:D:66:ARG:O	1:D:67:MET:C	2.58	0.42
1:D:53:LEU:HA	1:D:53:LEU:HD12	1.81	0.42
1:E:42:LEU:HD22	1:E:195:ALA:CB	2.50	0.42
1:E:97:LEU:HD21	1:E:129:LEU:HD21	2.02	0.41
1:E:36:THR:HA	1:E:223:ASN:O	2.20	0.41
1:C:53:LEU:HA	1:C:53:LEU:HD23	1.75	0.41
1:D:145:ILE:HD12	1:D:145:ILE:HG21	1.78	0.41
1:A:73:LYS:HZ3	2:Q:157:THR:HB	1.85	0.41
1:F:85:ILE:HD11	1:F:219:LEU:HD22	2.02	0.41
1:C:199:PHE:O	2:N:120:ASN:OD1	2.38	0.41
2:N:134:GLU:O	2:N:134:GLU:HG3	2.20	0.41
1:A:196:ALA:HA	1:A:197:PRO:HD2	1.90	0.41
1:B:97:LEU:HB3	1:B:191:TYR:CD2	2.56	0.41
1:F:39:LEU:CD1	1:F:224:PHE:CD1	2.99	0.41
1:C:82:GLY:HA3	1:C:125:MET:HE1	2.02	0.41
1:F:137:GLN:O	1:F:137:GLN:HG3	2.20	0.41
1:B:125:MET:HE3	1:B:177:VAL:HG21	2.03	0.41
1:F:52:PRO:CA	1:F:190:ASN:OD1	2.68	0.41
1:A:86:VAL:HG23	1:A:93:HIS:HB2	2.02	0.41
1:C:222:PHE:HB2	1:C:224:PHE:CZ	2.55	0.41
2:M:85:ILE:HD12	2:M:129:VAL:CG2	2.49	0.41
1:C:85:ILE:CD1	1:C:219:LEU:HD22	2.51	0.41
1:E:215:LEU:N	1:E:216:PRO:CD	2.84	0.41
1:B:131:ARG:HD2	1:B:135:VAL:HB	2.03	0.41
1:F:83:VAL:CG2	1:F:145:ILE:HD11	2.48	0.41
2:P:108:ASP:OD2	2:P:132:GLY:HA3	2.21	0.41
1:B:64:PHE:CD1	1:B:67:MET:CE	3.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:180:GLN:O	1:C:182:LYS:N	2.53	0.41
2:O:104:LEU:HD12	2:O:140:LEU:HD23	2.03	0.41
1:E:205:ALA:N	1:E:206:PRO:CD	2.84	0.41
1:C:101:THR:CG2	1:C:101:THR:O	2.66	0.41
1:F:50:LYS:HD3	1:F:186:ALA:O	2.21	0.41
1:C:89:HIS:O	1:C:90:ARG:HB2	2.21	0.41
1:F:35:ARG:NH2	1:F:143:ASP:OD2	2.54	0.41
1:D:35:ARG:NH1	1:D:35:ARG:CG	2.31	0.40
1:C:218:LEU:HD23	1:D:218:LEU:HD23	2.03	0.40
2:P:108:ASP:CG	2:P:108:ASP:O	2.59	0.40
1:A:145:ILE:HD12	1:A:145:ILE:HG21	1.85	0.40
1:F:73:LYS:C	1:F:74:ILE:HG22	2.40	0.40
1:B:84:LEU:HD12	1:B:177:VAL:HB	2.03	0.40
1:A:73:LYS:HD3	2:Q:157:THR:HG21	2.03	0.40
2:O:110:LEU:HA	2:O:110:LEU:HD23	1.73	0.40
1:D:121:LEU:HA	1:D:121:LEU:HD12	1.82	0.40
2:M:120:ASN:HA	2:M:120:ASN:HD22	1.49	0.40
1:C:79:THR:HB	1:C:172:LYS:HA	2.04	0.40
1:D:135:VAL:HG12	1:D:136:LEU:N	2.36	0.40
2:N:93:THR:O	2:N:96:ASP:HB2	2.21	0.40
1:A:205:ALA:N	1:A:206:PRO:HD2	2.37	0.40
1:F:37:ILE:HG23	1:F:37:ILE:HD13	1.87	0.40
1:D:84:LEU:HD12	1:D:84:LEU:N	2.37	0.40
1:A:91:LEU:HA	1:A:92:PRO:HD3	1.86	0.40

There are no symmetry-related clashes.

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	194/202 (96%)	188 (97%)	6 (3%)	0	100	100
1	B	196/202 (97%)	187 (95%)	8 (4%)	1 (0%)	38	70
1	C	194/202 (96%)	186 (96%)	7 (4%)	1 (0%)	38	70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	196/202 (97%)	184 (94%)	12 (6%)	0	100	100
1	E	187/202 (93%)	174 (93%)	12 (6%)	1 (0%)	38	70
1	F	195/202 (96%)	181 (93%)	13 (7%)	1 (0%)	38	70
2	L	72/90 (80%)	64 (89%)	7 (10%)	1 (1%)	16	41
2	M	79/90 (88%)	77 (98%)	1 (1%)	1 (1%)	18	43
2	N	78/90 (87%)	77 (99%)	1 (1%)	0	100	100
2	O	79/90 (88%)	75 (95%)	4 (5%)	0	100	100
2	P	77/90 (86%)	72 (94%)	4 (5%)	1 (1%)	18	43
2	Q	78/90 (87%)	70 (90%)	8 (10%)	0	100	100
All	All	1625/1752 (93%)	1535 (94%)	83 (5%)	7 (0%)	43	76

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	133	ASP
1	C	181	GLU
2	L	110	LEU
2	M	103	SER
2	P	103	SER
1	F	74	ILE
1	E	113	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	169/181 (93%)	159 (94%)	10 (6%)	28	56
1	B	172/181 (95%)	165 (96%)	7 (4%)	41	74
1	C	166/181 (92%)	159 (96%)	7 (4%)	40	73
1	D	173/181 (96%)	163 (94%)	10 (6%)	28	57
1	E	158/181 (87%)	147 (93%)	11 (7%)	21	47
1	F	168/181 (93%)	157 (94%)	11 (6%)	24	51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	L	43/78 (55%)	37 (86%)	6 (14%)	5	12
2	M	62/78 (80%)	57 (92%)	5 (8%)	17	36
2	N	63/78 (81%)	62 (98%)	1 (2%)	75	94
2	O	62/78 (80%)	59 (95%)	3 (5%)	35	68
2	P	59/78 (76%)	52 (88%)	7 (12%)	8	18
2	Q	57/78 (73%)	51 (90%)	6 (10%)	10	22
All	All	1352/1554 (87%)	1268 (94%)	84 (6%)	26	54

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LYS
1	A	57	ASP
1	A	63	ARG
1	A	101	THR
1	A	128	ILE
1	A	142	ASP
1	A	145	ILE
1	A	167	LYS
1	A	227	ASN
1	A	228	LEU
1	B	46	THR
1	B	74	ILE
1	B	101	THR
1	B	131	ARG
1	B	132	GLN
1	B	174	LEU
1	B	220	SER
1	C	132	GLN
1	C	133	ASP
1	C	137	GLN
1	C	144	CYS
1	C	180	GLN
1	C	190	ASN
1	C	194	VAL
1	D	35	ARG
1	D	46	THR
1	D	79	THR
1	D	99	LEU
1	D	101	THR

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Mol	Chain	Res	Type
1	D	102	THR
1	D	131	ARG
1	D	145	ILE
1	D	154	GLU
1	D	189	LYS
1	E	42	LEU
1	E	57	ASP
1	E	68	ARG
1	E	78	ARG
1	E	87	HIS
1	E	99	LEU
1	E	141	ILE
1	E	145	ILE
1	E	174	LEU
1	E	194	VAL
1	E	223	ASN
1	F	63	ARG
1	F	72	ASP
1	F	74	ILE
1	F	80	VAL
1	F	102	THR
1	F	116	ASP
1	F	131	ARG
1	F	132	GLN
1	F	135	VAL
1	F	137	GLN
1	F	178	GLN
2	L	109	ILE
2	L	112	ILE
2	L	118	ARG
2	L	128	LEU
2	L	131	VAL
2	L	133	SER
2	M	113	LYS
2	M	118	ARG
2	M	120	ASN
2	M	123	SER
2	M	131	VAL
2	N	83	LEU
2	O	131	VAL
2	O	157	THR
2	O	159	SER

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Mol	Chain	Res	Type
2	P	97	LEU
2	P	111	GLU
2	P	117	ASN
2	P	131	VAL
2	P	136	SER
2	P	152	GLN
2	P	157	THR
2	Q	109	ILE
2	Q	110	LEU
2	Q	112	ILE
2	Q	117	ASN
2	Q	131	VAL
2	Q	157	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	178	GLN
1	A	227	ASN
1	C	112	ASN
1	C	137	GLN
1	C	152	ASN
1	C	178	GLN
1	D	152	ASN
1	E	223	ASN
1	E	227	ASN
1	F	178	GLN
2	L	153	ASN
2	M	120	ASN
2	M	122	GLN
2	N	102	HIS
2	N	152	GLN
2	O	102	HIS
2	O	120	ASN
2	O	160	ASN
2	P	102	HIS
2	P	117	ASN
2	P	152	GLN
2	Q	102	HIS
2	Q	117	ASN
2	Q	153	ASN

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 ⓘ

There are no ligands in this entry.

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	196/202 (97%)	-0.33	4 (2%) 62 68	14, 31, 48, 59	0
1	B	198/202 (98%)	-0.25	1 (0%) 88 92	16, 29, 54, 60	0
1	C	196/202 (97%)	0.05	10 (5%) 27 30	15, 42, 69, 77	0
1	D	198/202 (98%)	-0.31	6 (3%) 48 54	13, 24, 52, 61	0
1	E	191/202 (94%)	-0.07	4 (2%) 60 67	24, 43, 76, 81	0
1	F	197/202 (97%)	-0.15	6 (3%) 48 54	23, 39, 70, 76	0
2	L	76/90 (84%)	0.50	4 (5%) 25 28	38, 63, 95, 98	0
2	M	81/90 (90%)	-0.22	1 (1%) 75 81	18, 33, 45, 47	0
2	N	80/90 (88%)	-0.38	0 100 100	18, 31, 37, 38	0
2	O	81/90 (90%)	-0.17	0 100 100	22, 36, 49, 51	0
2	P	79/90 (87%)	-0.15	0 100 100	26, 40, 56, 58	0
2	Q	80/90 (88%)	0.15	2 (2%) 54 61	36, 54, 66, 67	0
All	All	1653/1752 (94%)	-0.14	38 (2%) 57 64	13, 36, 67, 98	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	105	GLY	5.4
1	F	59	SER	4.8
1	D	60	VAL	4.7
1	C	54	TYR	4.4
1	B	54	TYR	3.6
1	D	54	TYR	3.6
1	F	58	SER	3.1
2	Q	110	LEU	3.0
2	L	102	HIS	2.9
1	C	101	THR	2.9
2	Q	152	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	34	GLU	2.8
1	E	54	TYR	2.8
1	E	228	LEU	2.7
1	F	60	VAL	2.7
1	C	102	THR	2.6
1	C	136	LEU	2.6
1	D	61	ALA	2.6
2	M	133	SER	2.6
1	C	55	GLU	2.6
1	C	69	GLU	2.5
2	L	106	VAL	2.5
1	D	231	HIS	2.4
1	A	134	GLY	2.3
1	A	133	ASP	2.3
1	F	54	TYR	2.3
1	F	57	ASP	2.3
1	D	58	SER	2.3
1	C	228	LEU	2.3
1	A	136	LEU	2.2
1	A	135	VAL	2.2
1	E	72	ASP	2.2
1	C	56	LYS	2.2
2	L	118	ARG	2.1
1	E	113	PRO	2.1
1	C	74	ILE	2.1
1	C	229	GLU	2.0
1	F	102	THR	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.