



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 9, 2018 – 10:59 PM EST

PDB ID : 4UW8
Title : Structure of the carboxy-terminal domain of the bacteriophage T5 L- shaped tail fiber with its intra-molecular chaperone domain
Authors : Garcia-Doval, C.; Luque, D.; Caston, J.R.; Otero, J.M.; Llamas-Saiz, A.L.; Boulanger, P.; van Raaij, M.J.
Deposited on : 2014-08-08
Resolution : 2.52 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030736

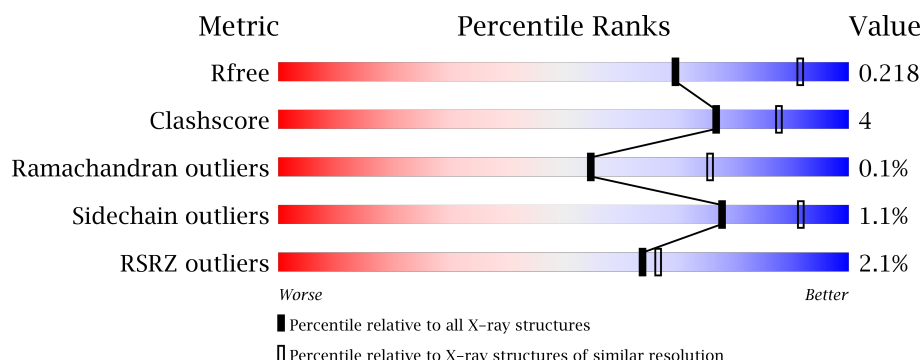
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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}	100719	4636 (2.54-2.50)
Clashscore	112137	5382 (2.54-2.50)
Ramachandran outliers	110173	5282 (2.54-2.50)
Sidechain outliers	110143	5284 (2.54-2.50)
RSRZ outliers	101464	4669 (2.54-2.50)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	427	<div> <div>0.1%</div> <div> <div></div> <div>87%</div> <div>7%</div> <div>6%</div> </div> </div>
1	B	427	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>6%</div> <div>6%</div> </div> </div>
1	C	427	<div> <div>0.1%</div> <div> <div></div> <div>87%</div> <div>7%</div> <div>6%</div> </div> </div>
1	D	427	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>5%</div> <div>6%</div> </div> </div>
1	E	427	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>6%</div> <div>6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	427	 2% 86% 7% • 6%
1	G	427	 2% 87% 6% • 6%
1	H	427	 4% 87% 6% • 6%
1	I	427	 % 88% 5% • 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FLC	A	2397	-	-	-	X
2	FLC	B	2397	-	-	-	X
2	FLC	C	2397	-	-	X	X
2	FLC	E	2397	-	-	-	X
2	FLC	F	2397	-	-	-	X
2	FLC	G	2397	-	-	-	X
2	FLC	I	2397	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 28644 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called L-SHAPED TAIL FIBER PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	B	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	C	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	D	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	E	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	F	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			
1	G	400	Total	C	N	O	S	0	2	0
			3058	1919	541	585	13			
1	H	401	Total	C	N	O	S	0	3	0
			3071	1928	542	588	13			
1	I	401	Total	C	N	O	S	0	2	0
			3066	1925	542	586	13			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1264	ALA	SER	engineered mutation	UNP P13390
B	1264	ALA	SER	engineered mutation	UNP P13390
C	1264	ALA	SER	engineered mutation	UNP P13390
D	1264	ALA	SER	engineered mutation	UNP P13390
E	1264	ALA	SER	engineered mutation	UNP P13390
F	1264	ALA	SER	engineered mutation	UNP P13390
G	1264	ALA	SER	engineered mutation	UNP P13390
H	1264	ALA	SER	engineered mutation	UNP P13390
I	1264	ALA	SER	engineered mutation	UNP P13390

- Molecule 2 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		
2	E	1	Total	C	O	0	0
			13	6	7		
2	F	1	Total	C	O	0	0
			13	6	7		
2	G	1	Total	C	O	0	0
			13	6	7		
2	H	1	Total	C	O	0	0
			13	6	7		
2	I	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	154	Total	O	0	0
			154	154		
3	B	128	Total	O	0	0
			128	128		
3	C	108	Total	O	0	0
			108	108		

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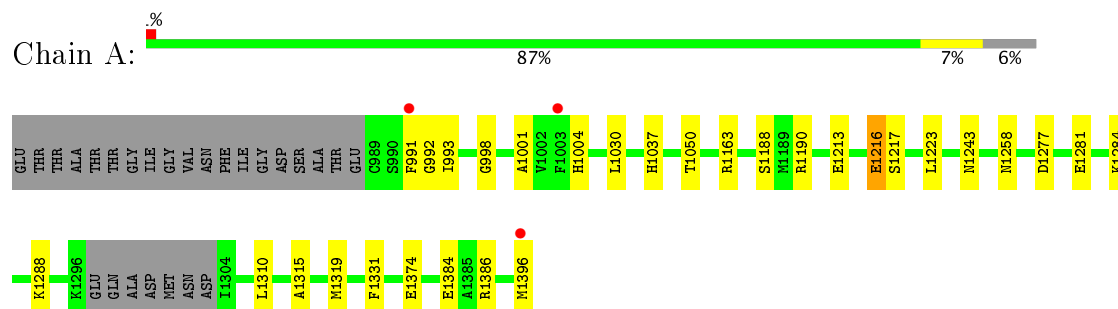
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	136	Total 136	O 136	0	0
3	E	122	Total 122	O 122	0	0
3	F	65	Total 65	O 65	0	0
3	G	90	Total 90	O 90	0	0
3	H	63	Total 63	O 63	0	0
3	I	70	Total 70	O 70	0	0

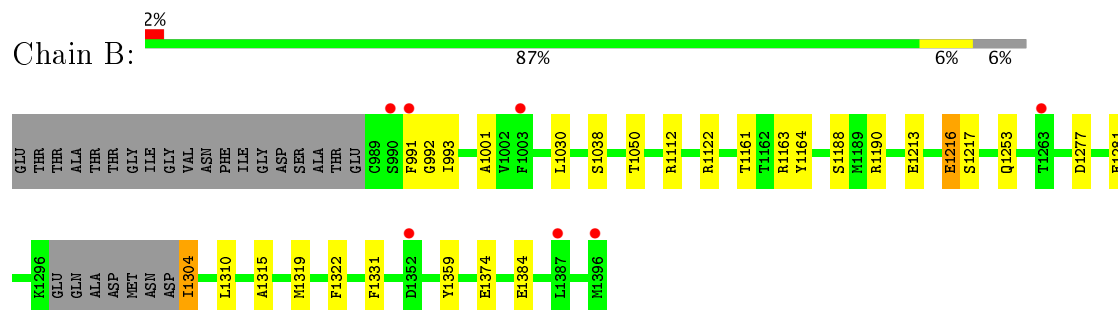
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

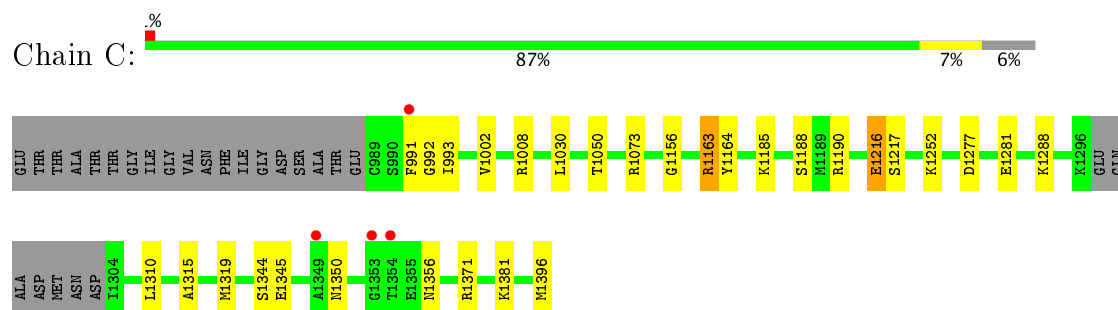
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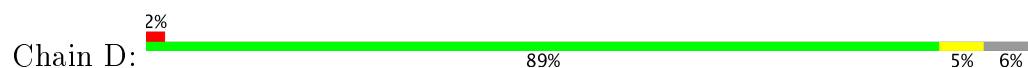
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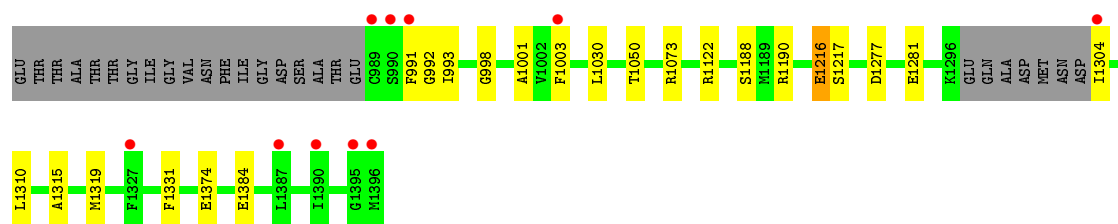


• Molecule 1: L-SHAPED TAIL FIBER PROTEIN

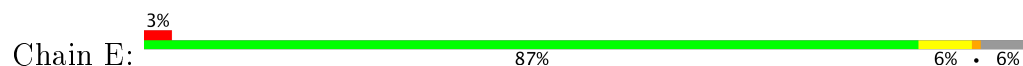


• Molecule 1: L-SHAPED TAIL FIBER PROTEIN

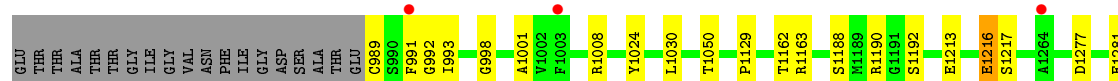
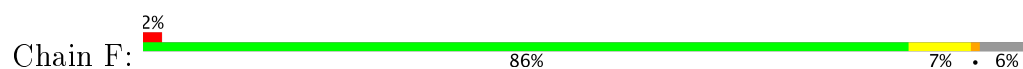




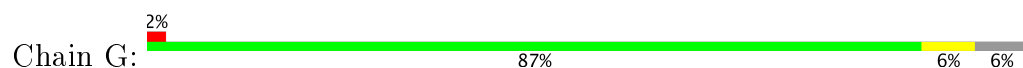
• Molecule 1: L-SHAPED TAIL FIBER PROTEIN



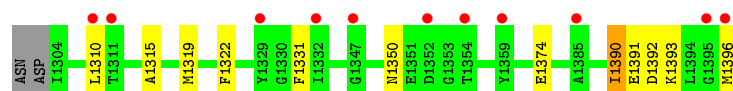
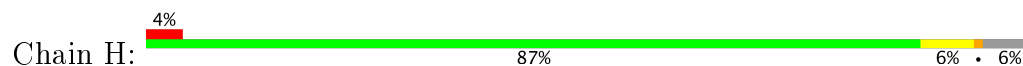
• Molecule 1: L-SHAPED TAIL FIBER PROTEIN



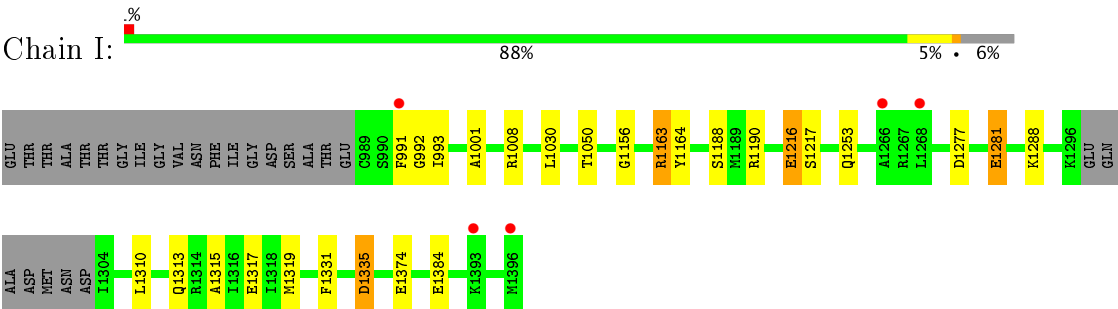
• Molecule 1: L-SHAPED TAIL FIBER PROTEIN



• Molecule 1: L-SHAPED TAIL FIBER PROTEIN



● Molecule 1: L-SHAPED TAIL FIBER PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	160.85Å 99.29Å 286.23Å 90.00° 91.51° 90.00°	Depositor
Resolution (Å)	95.38 – 2.52 95.38 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.5 (95.38-2.52) 98.5 (95.38-2.52)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.185 , 0.213 0.190 , 0.218	Depositor DCC
R_{free} test set	2061 reflections (1.39%)	DCC
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	28644	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

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

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.55	2/3147 (0.1%)	0.72	1/4269 (0.0%)
1	B	0.57	2/3147 (0.1%)	0.74	1/4269 (0.0%)
1	C	0.56	0/3147	0.73	0/4269
1	D	0.52	0/3147	0.72	0/4269
1	E	0.56	2/3147 (0.1%)	0.74	1/4269 (0.0%)
1	F	0.55	0/3147	0.76	4/4269 (0.1%)
1	G	0.50	0/3139	0.76	3/4258 (0.1%)
1	H	0.50	0/3155	0.71	0/4280
1	I	0.54	1/3147 (0.0%)	0.73	3/4269 (0.1%)
All	All	0.54	7/28323 (0.0%)	0.73	13/38421 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1213	GLU	CD-OE1	-5.92	1.19	1.25
1	B	1213	GLU	CD-OE1	-5.86	1.19	1.25
1	E	1355	GLU	CD-OE2	-5.42	1.19	1.25
1	A	1213	GLU	CD-OE2	-5.24	1.19	1.25
1	B	1213	GLU	CD-OE2	-5.21	1.20	1.25
1	I	1281	GLU	CD-OE1	5.21	1.31	1.25
1	E	1335	ASP	CB-CG	5.00	1.62	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	1335	ASP	CB-CG-OD2	9.50	126.85	118.30
1	I	1335	ASP	CB-CG-OD2	9.34	126.70	118.30
1	G	1122[A]	ARG	CG-CD-NE	8.76	130.19	111.80
1	G	1122[B]	ARG	CG-CD-NE	8.76	130.19	111.80
1	E	1335	ASP	CB-CG-OD2	8.70	126.13	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	1281	GLU	CG-CD-OE2	-7.23	103.84	118.30
1	F	1384	GLU	OE1-CD-OE2	-6.93	114.98	123.30
1	I	1281	GLU	CG-CD-OE1	6.79	131.87	118.30
1	G	1386	ARG	CG-CD-NE	6.66	125.79	111.80
1	F	1319	MET	CG-SD-CE	-5.98	90.64	100.20
1	F	1384	GLU	CG-CD-OE2	5.15	128.60	118.30
1	B	1163	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	A	1163	ARG	NE-CZ-NH2	-5.04	117.78	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3066	0	2963	26	0
1	B	3066	0	2963	29	0
1	C	3066	0	2963	29	0
1	D	3066	0	2963	24	0
1	E	3066	0	2963	25	0
1	F	3066	0	2963	32	0
1	G	3058	0	2952	26	0
1	H	3071	0	2967	25	0
1	I	3066	0	2963	23	0
2	A	13	0	5	2	0
2	B	13	0	5	2	0
2	C	13	0	5	4	0
2	D	13	0	5	2	0
2	E	13	0	5	0	0
2	F	13	0	5	1	0
2	G	13	0	5	0	0
2	H	13	0	5	1	0
2	I	13	0	5	0	0
3	A	154	0	0	0	0
3	B	128	0	0	5	0
3	C	108	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	136	0	0	2	0
3	E	122	0	0	1	0
3	F	65	0	0	3	0
3	G	90	0	0	3	0
3	H	63	0	0	2	0
3	I	70	0	0	0	0
All	All	28644	0	26705	191	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (191) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1331:PHE:HB3	1:B:1374:GLU:HG2	1.28	1.15
1:F:989:CYS:SG	3:F:2002:HOH:O	2.25	0.92
1:D:1304:ILE:N	3:D:2119:HOH:O	2.08	0.86
1:H:1281:GLU:HB3	1:H:1322:PHE:CE2	2.16	0.80
1:B:1281:GLU:HB3	1:B:1322:PHE:CE2	2.17	0.79
1:F:1335:ASP:OD1	3:F:2054:HOH:O	2.00	0.78
1:G:1335:ASP:OD2	3:G:2080:HOH:O	2.02	0.77
1:C:1073:ARG:HH22	2:C:2397:FLC:CA	1.97	0.76
1:D:993:ILE:HB	1:E:991[B]:PHE:CE1	2.21	0.75
1:B:1164:TYR:CE2	1:E:1112:ARG:HD2	2.21	0.75
1:B:993:ILE:HB	1:C:991[A]:PHE:CE1	2.21	0.74
1:G:991[B]:PHE:CE1	1:I:993:ILE:HB	2.27	0.69
1:F:1313:GLN:O	1:F:1317:GLU:HG2	1.93	0.68
1:D:993:ILE:HD12	1:E:991[B]:PHE:CD1	2.28	0.68
1:G:1394:LEU:HD12	1:G:1396:MET:HE1	1.75	0.68
1:A:993:ILE:HB	1:B:991[A]:PHE:CE1	2.29	0.67
1:I:1313:GLN:O	1:I:1317:GLU:HG2	1.94	0.67
1:F:1162:THR:HG22	3:F:2036:HOH:O	1.95	0.66
1:C:1073:ARG:HH22	2:C:2397:FLC:HA2	1.64	0.63
1:G:1316:ILE:HG22	1:G:1320:GLU:OE2	1.98	0.63
1:F:1192:SER:OG	1:F:1213:GLU:OE1	2.18	0.62
1:G:1316:ILE:CG2	1:G:1320:GLU:OE2	2.48	0.62
1:B:993:ILE:HD12	1:C:991[A]:PHE:CD1	2.36	0.61
1:H:993:ILE:HB	1:I:991[B]:PHE:CE1	2.36	0.60
1:H:1160[A]:ASP:OD1	1:H:1162:THR:O	2.22	0.58
3:B:2056:HOH:O	1:C:1163:ARG:HD2	2.03	0.57
2:F:2397:FLC:OG2	2:F:2397:FLC:CBC	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1185:LYS:HE2	3:C:2065:HOH:O	2.04	0.57
1:G:993:ILE:HB	1:H:991[B]:PHE:CE1	2.39	0.57
1:E:1288:LYS:HG2	1:F:1384:GLU:OE2	2.04	0.56
1:E:993:ILE:HB	1:F:991[B]:PHE:CE1	2.41	0.56
1:H:1390:ILE:O	1:H:1393:LYS:N	2.38	0.55
1:H:993:ILE:HD12	1:I:991[B]:PHE:CD1	2.42	0.55
1:C:1073:ARG:HH22	2:C:2397:FLC:HA1	1.70	0.54
1:B:1161:THR:HG23	3:B:2069:HOH:O	2.07	0.54
2:D:2397:FLC:OA2	2:D:2397:FLC:OHB	2.19	0.54
1:G:1118:ILE:O	1:G:1122[A]:ARG:NH1	2.41	0.54
1:A:991[A]:PHE:CE1	1:C:993:ILE:HB	2.43	0.53
1:C:1156:GLY:HA2	1:C:1164:TYR:CE2	2.43	0.53
1:H:1156:GLY:HA2	1:H:1164:TYR:CE2	2.44	0.53
1:D:993:ILE:HB	1:E:991[B]:PHE:CZ	2.43	0.53
1:G:1394:LEU:HD12	1:G:1396:MET:CE	2.37	0.53
3:H:2026:HOH:O	1:I:1163:ARG:HD2	2.08	0.53
1:F:1394:LEU:HD12	1:F:1396:MET:HE1	1.90	0.52
1:I:1315:ALA:O	1:I:1319:MET:HG2	2.10	0.52
1:A:1216:GLU:HG3	1:A:1217:SER:N	2.25	0.52
1:B:1112:ARG:HD3	3:B:2058:HOH:O	2.09	0.52
1:B:1315:ALA:O	1:B:1319:MET:HG2	2.10	0.52
1:E:1156:GLY:HA2	1:E:1164:TYR:CE2	2.45	0.52
1:G:1384:GLU:OE2	1:I:1288:LYS:HE2	2.10	0.52
1:D:1315:ALA:O	1:D:1319:MET:HG2	2.10	0.51
3:D:2063:HOH:O	1:E:1163:ARG:HD2	2.10	0.51
1:E:1315:ALA:O	1:E:1319:MET:HG2	2.10	0.51
1:B:993:ILE:HB	1:C:991[A]:PHE:CZ	2.45	0.51
1:G:1216:GLU:HG3	1:G:1217:SER:N	2.25	0.51
1:H:1315:ALA:O	1:H:1319:MET:HG2	2.10	0.51
1:B:1277:ASP:O	1:B:1281:GLU:HG2	2.11	0.51
1:C:1315:ALA:O	1:C:1319:MET:HG2	2.11	0.51
1:A:991[A]:PHE:CD1	1:C:993:ILE:HD12	2.46	0.51
1:E:1277:ASP:O	1:E:1281:GLU:HG2	2.11	0.51
1:A:1277:ASP:O	1:A:1281:GLU:HG2	2.11	0.51
1:B:1216:GLU:HG3	1:B:1217:SER:N	2.25	0.51
1:I:1156:GLY:HA2	1:I:1164:TYR:CE2	2.45	0.51
1:B:1359:TYR:HD2	3:B:2123:HOH:O	1.94	0.51
1:A:993:ILE:HD12	1:B:991[A]:PHE:CD1	2.45	0.51
1:G:1315:ALA:O	1:G:1319:MET:HG2	2.10	0.51
1:H:1277:ASP:O	1:H:1281:GLU:HG2	2.11	0.51
1:H:1281:GLU:HB3	1:H:1322:PHE:HE2	1.72	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1277:ASP:O	1:C:1281:GLU:HG2	2.11	0.50
1:F:1277:ASP:O	1:F:1281:GLU:HG2	2.11	0.50
1:G:1277:ASP:O	1:G:1281:GLU:HG2	2.11	0.50
1:A:1223:LEU:HD23	1:C:1345:GLU:HA	1.92	0.50
1:D:1277:ASP:O	1:D:1281:GLU:HG2	2.11	0.50
1:F:1216:GLU:HG3	1:F:1217:SER:N	2.25	0.50
1:H:1350:ASN:HB2	3:H:2059:HOH:O	2.12	0.50
1:E:1216:GLU:HG3	1:E:1217:SER:N	2.26	0.50
1:C:1216:GLU:HG3	1:C:1217:SER:N	2.27	0.50
1:A:1315:ALA:O	1:A:1319:MET:HG2	2.11	0.50
1:A:1030:LEU:HD23	1:B:1050:THR:HG22	1.92	0.50
1:E:1305:ARG:NH2	1:F:1332:ILE:O	2.37	0.49
1:F:1394:LEU:HB2	1:F:1396:MET:HE2	1.94	0.49
1:I:1216:GLU:HG3	1:I:1217:SER:N	2.28	0.49
1:D:991[B]:PHE:CE1	1:F:993:ILE:HB	2.47	0.49
1:D:998:GLY:HA2	1:F:1008:ARG:O	2.12	0.49
1:D:1216:GLU:HG3	1:D:1217:SER:N	2.27	0.48
1:A:1037:HIS:ND1	2:A:2397:FLC:OB2	2.32	0.48
3:B:2102:HOH:O	1:C:1381:LYS:HE2	2.12	0.48
1:A:1050:THR:HG22	1:C:1030:LEU:HD23	1.95	0.48
1:G:1119:GLN:HA	1:G:1122[A]:ARG:CD	2.43	0.48
1:F:1394:LEU:HD12	1:F:1396:MET:CE	2.43	0.47
1:H:1216:GLU:HG3	1:H:1217:SER:N	2.29	0.47
1:D:1122[B]:ARG:NH1	1:F:1129:PRO:O	2.47	0.47
1:E:993:ILE:HD12	1:F:991[B]:PHE:CD1	2.49	0.47
1:D:1030:LEU:HD23	1:E:1050:THR:HG22	1.97	0.47
1:A:1223:LEU:HD22	1:C:1344:SER:O	2.15	0.47
1:B:1281:GLU:HB3	1:B:1322:PHE:HE2	1.74	0.47
1:A:1004:HIS:NE2	2:C:2397:FLC:OG2	2.44	0.47
1:D:1073:ARG:NH2	2:D:2397:FLC:OG1	2.47	0.47
1:A:1384:GLU:OE2	1:C:1288:LYS:HE2	2.14	0.46
1:A:991[B]:PHE:CD1	1:A:992:GLY:N	2.84	0.46
1:A:1288:LYS:HE2	1:B:1384:GLU:OE2	2.16	0.46
1:H:1390:ILE:O	1:H:1392:ASP:N	2.49	0.46
1:H:1030:LEU:HD23	1:I:1050:THR:HG22	1.98	0.46
1:D:993:ILE:HD12	1:E:991[B]:PHE:HD1	1.79	0.46
1:G:1050:THR:HG22	1:I:1030:LEU:HD23	1.98	0.46
1:G:991[B]:PHE:CD1	1:I:993:ILE:HD12	2.50	0.46
3:G:2035:HOH:O	1:H:1163:ARG:HD2	2.15	0.46
1:B:1281:GLU:CB	1:B:1322:PHE:CE2	2.97	0.46
1:E:991[A]:PHE:CD1	1:E:992:GLY:N	2.84	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:991[B]:PHE:CD1	1:C:992:GLY:N	2.84	0.46
1:I:1277:ASP:O	1:I:1281:GLU:HG2	2.16	0.46
1:B:991[B]:PHE:CD1	1:B:992:GLY:N	2.85	0.45
2:B:2397:FLC:HG1	1:C:1002:VAL:HG11	1.99	0.45
1:E:1030:LEU:HD23	1:F:1050:THR:HG22	1.98	0.45
1:C:1350:ASN:OD1	1:C:1356:ASN:OD1	2.35	0.45
1:I:991[A]:PHE:CD1	1:I:992:GLY:N	2.85	0.45
1:F:1162:THR:O	1:F:1162:THR:HG23	2.17	0.45
1:H:991[A]:PHE:CD1	1:H:992:GLY:N	2.85	0.44
2:A:2397:FLC:OB2	2:A:2397:FLC:CGC	2.63	0.44
1:D:991[A]:PHE:CD1	1:D:992:GLY:N	2.85	0.44
1:B:1188:SER:OG	1:B:1190:ARG:NH1	2.51	0.44
1:D:1050:THR:HG22	1:F:1030:LEU:HD23	1.98	0.44
1:H:1188:SER:OG	1:H:1190:ARG:NH1	2.50	0.44
1:G:991[B]:PHE:CZ	1:I:993:ILE:HB	2.53	0.44
1:A:1188:SER:OG	1:A:1190:ARG:NH1	2.51	0.44
1:A:998:GLY:HA2	1:C:1008:ARG:O	2.17	0.44
1:F:1188:SER:OG	1:F:1190:ARG:NH1	2.50	0.43
1:G:1394:LEU:HB2	1:G:1396:MET:HE2	1.99	0.43
1:B:1030:LEU:HD23	1:C:1050:THR:HG22	2.00	0.43
1:H:1160[A]:ASP:OD2	1:H:1164:TYR:HE1	2.01	0.43
1:D:991[B]:PHE:CZ	1:F:993:ILE:HB	2.53	0.43
1:F:991[A]:PHE:CD1	1:F:992:GLY:N	2.86	0.43
1:G:991[A]:PHE:CD1	1:G:992:GLY:N	2.86	0.43
1:I:1188:SER:OG	1:I:1190:ARG:NH1	2.51	0.43
1:E:1008:ARG:O	1:F:998:GLY:HA2	2.19	0.42
2:H:2397:FLC:OHB	2:H:2397:FLC:OA2	2.32	0.42
1:C:1188:SER:OG	1:C:1190:ARG:NH1	2.51	0.42
1:H:1281:GLU:CB	1:H:1322:PHE:CE2	2.95	0.42
1:H:1331:PHE:HB3	1:H:1374:GLU:HB3	2.02	0.42
1:E:1188:SER:OG	1:E:1190:ARG:NH1	2.52	0.42
1:A:1243:ASN:OD1	1:C:1252:LYS:HB3	2.20	0.42
1:B:993:ILE:HG13	1:B:1001:ALA:HB1	2.01	0.42
1:E:1334:TYR:HE2	1:E:1336:LYS:HE3	1.84	0.42
1:G:1030:LEU:HD23	1:H:1050:THR:HG22	2.02	0.42
1:E:993:ILE:HG13	1:E:1001:ALA:HB1	2.02	0.42
1:G:1188:SER:OG	1:G:1190:ARG:NH1	2.52	0.42
1:G:1371:ARG:NH1	1:I:1253:GLN:OE1	2.42	0.42
1:A:993:ILE:HG13	1:A:1001:ALA:HB1	2.02	0.41
1:E:1331:PHE:HB3	1:E:1374:GLU:HB3	2.02	0.41
1:F:993:ILE:HG13	1:F:1001:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:993:ILE:HG13	1:D:1001:ALA:HB1	2.01	0.41
1:I:1331:PHE:HB3	1:I:1374:GLU:HB3	2.02	0.41
1:D:991[B]:PHE:CD1	1:F:993:ILE:HD12	2.56	0.41
1:B:1304:ILE:HG13	1:B:1304:ILE:O	2.20	0.41
1:G:1163:ARG:HD2	3:G:2048:HOH:O	2.20	0.41
1:I:993:ILE:HG13	1:I:1001:ALA:HB1	2.01	0.41
1:D:1003:PHE:CD1	1:F:1024:TYR:CE2	3.09	0.41
1:D:1384:GLU:OE2	1:F:1288:LYS:HE2	2.20	0.41
1:G:1331:PHE:HB3	1:G:1374:GLU:HB3	2.02	0.41
1:H:993:ILE:HG13	1:H:1001:ALA:HB1	2.03	0.41
1:A:1223:LEU:CD2	1:C:1345:GLU:HB3	2.50	0.41
1:E:1351:GLU:HG2	3:E:2110:HOH:O	2.20	0.41
1:A:1331:PHE:HB3	1:A:1374:GLU:HB3	2.03	0.41
1:G:998:GLY:HA2	1:I:1008:ARG:O	2.21	0.41
1:A:1284:LYS:HD3	1:A:1386:ARG:CG	2.51	0.41
1:A:1258:ASN:HB2	1:B:1253:GLN:HB3	2.03	0.41
1:D:1331:PHE:HB3	1:D:1374:GLU:HB3	2.02	0.41
1:B:1122[B]:ARG:HD3	1:B:1122[B]:ARG:HA	1.80	0.40
1:B:1253:GLN:OE1	1:C:1371:ARG:NH1	2.47	0.40
1:B:1038:SER:HB2	2:B:2397:FLC:OB2	2.21	0.40
1:D:1188:SER:OG	1:D:1190:ARG:NH1	2.53	0.40
1:F:1331:PHE:HB3	1:F:1374:GLU:HB3	2.04	0.40
1:F:1304:ILE:O	1:F:1304:ILE:HG22	2.22	0.40
1:H:1288:LYS:HE2	1:I:1384:GLU:OE2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	399/427 (93%)	390 (98%)	9 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	399/427 (93%)	389 (98%)	10 (2%)	0	100	100
1	C	399/427 (93%)	390 (98%)	9 (2%)	0	100	100
1	D	399/427 (93%)	390 (98%)	9 (2%)	0	100	100
1	E	399/427 (93%)	390 (98%)	9 (2%)	0	100	100
1	F	399/427 (93%)	390 (98%)	9 (2%)	0	100	100
1	G	398/427 (93%)	389 (98%)	9 (2%)	0	100	100
1	H	400/427 (94%)	388 (97%)	10 (2%)	2 (0%)	32	52
1	I	399/427 (93%)	390 (98%)	9 (2%)	0	100	100
All	All	3591/3843 (93%)	3506 (98%)	83 (2%)	2 (0%)	55	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	1390	ILE
1	H	1391	GLU

5.3.2 Protein sidechains [i](#)

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	318/336 (95%)	315 (99%)	3 (1%)	82	93
1	B	318/336 (95%)	315 (99%)	3 (1%)	82	93
1	C	318/336 (95%)	314 (99%)	4 (1%)	73	90
1	D	318/336 (95%)	316 (99%)	2 (1%)	89	96
1	E	318/336 (95%)	313 (98%)	5 (2%)	68	87
1	F	318/336 (95%)	313 (98%)	5 (2%)	68	87
1	G	317/336 (94%)	315 (99%)	2 (1%)	89	96
1	H	319/336 (95%)	315 (99%)	4 (1%)	73	90
1	I	318/336 (95%)	314 (99%)	4 (1%)	73	90
All	All	2862/3024 (95%)	2830 (99%)	32 (1%)	78	91

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1216	GLU
1	A	1310	LEU
1	A	1396	MET
1	B	1216	GLU
1	B	1304	ILE
1	B	1310	LEU
1	C	1163	ARG
1	C	1216	GLU
1	C	1310	LEU
1	C	1396	MET
1	D	1216	GLU
1	D	1310	LEU
1	E	1163	ARG
1	E	1216	GLU
1	E	1310	LEU
1	E	1335	ASP
1	E	1396	MET
1	F	1163	ARG
1	F	1216	GLU
1	F	1310	LEU
1	F	1319	MET
1	F	1335	ASP
1	G	1216	GLU
1	G	1310	LEU
1	H	1163	ARG
1	H	1216	GLU
1	H	1310	LEU
1	H	1396	MET
1	I	1163	ARG
1	I	1216	GLU
1	I	1310	LEU
1	I	1335	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FLC	A	2397	-	3,12,12	0.43	0	3,17,17	0.74	0
2	FLC	B	2397	-	3,12,12	0.65	0	3,17,17	0.80	0
2	FLC	C	2397	-	3,12,12	1.40	1 (33%)	3,17,17	2.06	1 (33%)
2	FLC	D	2397	-	3,12,12	0.62	0	3,17,17	1.57	1 (33%)
2	FLC	E	2397	-	3,12,12	0.72	0	3,17,17	0.69	0
2	FLC	F	2397	-	3,12,12	1.14	0	3,17,17	1.88	1 (33%)
2	FLC	G	2397	-	3,12,12	0.32	0	3,17,17	0.40	0
2	FLC	H	2397	-	3,12,12	0.80	0	3,17,17	1.71	1 (33%)
2	FLC	I	2397	-	3,12,12	0.32	0	3,17,17	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FLC	A	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	B	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	C	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	D	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	E	2397	-	-	0/6/16/16	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FLC	F	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	G	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	H	2397	-	-	0/6/16/16	0/0/0/0
2	FLC	I	2397	-	-	0/6/16/16	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2397	FLC	CA-CB	-2.23	1.51	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2397	FLC	CB-CG-CGC	-3.42	109.61	114.95
2	F	2397	FLC	CB-CG-CGC	-3.03	110.22	114.95
2	H	2397	FLC	CB-CA-CAC	-2.71	110.72	114.95
2	D	2397	FLC	CB-CG-CGC	-2.25	111.44	114.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2397	FLC	2	0
2	B	2397	FLC	2	0
2	C	2397	FLC	4	0
2	D	2397	FLC	2	0
2	F	2397	FLC	1	0
2	H	2397	FLC	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	401/427 (93%)	0.11	3 (0%) 87 88	19, 34, 68, 91	0
1	B	401/427 (93%)	0.09	7 (1%) 70 73	20, 34, 73, 93	0
1	C	401/427 (93%)	0.06	4 (0%) 82 84	20, 33, 68, 94	0
1	D	401/427 (93%)	0.10	10 (2%) 58 61	21, 36, 83, 115	0
1	E	401/427 (93%)	0.23	12 (2%) 51 54	20, 36, 86, 111	0
1	F	401/427 (93%)	0.12	9 (2%) 62 65	21, 36, 72, 106	0
1	G	400/427 (93%)	0.20	10 (2%) 58 61	25, 40, 85, 105	0
1	H	401/427 (93%)	0.32	15 (3%) 42 45	24, 42, 90, 120	0
1	I	401/427 (93%)	0.12	5 (1%) 79 81	27, 40, 74, 121	0
All	All	3608/3843 (93%)	0.15	75 (2%) 64 66	19, 37, 82, 121	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	991[A]	PHE	5.7
1	I	1396	MET	5.1
1	D	1396	MET	4.6
1	F	1396	MET	4.5
1	D	991[A]	PHE	4.3
1	F	1264	ALA	4.3
1	E	1282	ALA	4.2
1	H	1354	THR	4.1
1	A	991[A]	PHE	4.1
1	E	1359	TYR	4.0
1	I	991[A]	PHE	3.9
1	E	991[A]	PHE	3.9
1	G	1329	TYR	3.7
1	E	1333	CYS	3.7
1	H	1395	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	991[A]	PHE	3.5
1	H	1396	MET	3.5
1	B	1003	PHE	3.4
1	E	1385	ALA	3.4
1	H	1281	GLU	3.4
1	C	991[A]	PHE	3.3
1	E	1349	ALA	3.2
1	F	991[A]	PHE	3.2
1	A	1003	PHE	3.1
1	E	1353	GLY	3.1
1	H	1329	TYR	3.1
1	I	1268	LEU	3.0
1	E	1396	MET	3.0
1	G	1327	PHE	2.9
1	H	1292	PHE	2.8
1	C	1353	GLY	2.8
1	H	1310	LEU	2.8
1	G	1316	ILE	2.7
1	H	1311	THR	2.7
1	G	991[A]	PHE	2.7
1	H	1385	ALA	2.7
1	G	1295	TRP	2.7
1	E	1383	PHE	2.6
1	D	990	SER	2.6
1	B	990	SER	2.6
1	B	1396	MET	2.5
1	F	1389	ALA	2.5
1	I	1266	ALA	2.5
1	H	1347	GLY	2.5
1	E	1286	ILE	2.4
1	G	1319	MET	2.4
1	D	1395	GLY	2.4
1	D	989	CYS	2.4
1	H	989	CYS	2.4
1	D	1387	LEU	2.4
1	G	1331	PHE	2.4
1	C	1349	ALA	2.4
1	C	1354	THR	2.4
1	F	1387	LEU	2.4
1	D	1327	PHE	2.3
1	H	1352	ASP	2.3
1	I	1393	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	1359	TYR	2.3
1	G	1335	ASP	2.3
1	A	1396	MET	2.3
1	B	1387	LEU	2.3
1	F	1373	GLU	2.2
1	E	1368	TYR	2.2
1	G	1368	TYR	2.2
1	D	1304	ILE	2.2
1	B	1263	THR	2.2
1	E	1370	PHE	2.2
1	B	1352	ASP	2.2
1	F	1391	GLU	2.2
1	F	1388	SER	2.1
1	H	1332	ILE	2.1
1	F	1003	PHE	2.1
1	D	1003	PHE	2.1
1	D	1390	ILE	2.0
1	G	1306	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	FLC	I	2397	13/13	0.89	0.31	4.17	69,75,88,90	0
2	FLC	G	2397	13/13	0.94	0.32	3.90	71,76,96,98	0
2	FLC	F	2397	13/13	0.91	0.32	3.67	58,78,90,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FLC	B	2397	13/13	0.92	0.26	3.41	62,74,87,87	0
2	FLC	A	2397	13/13	0.94	0.37	2.85	80,87,97,101	0
2	FLC	E	2397	13/13	0.94	0.26	2.62	53,72,81,85	0
2	FLC	C	2397	13/13	0.90	0.28	2.24	66,79,90,90	0
2	FLC	H	2397	13/13	0.91	0.25	1.91	74,88,92,94	0
2	FLC	D	2397	13/13	0.94	0.19	0.84	61,79,88,88	0

6.5 Other polymers [i](#)

There are no such residues in this entry.