



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 11, 2016 – 05:03 PM EDT

PDB ID : 5FCL
Title : Crystal structure of Cas1 from *Pectobacterium atrosepticum*
Authors : Wilkinson, M.E.; Nakatani, Y.; Opel-Reading, H.K.; Fineran, P.C.; Krause, K.L.
Deposited on : 2015-12-15
Resolution : 2.70 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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-20027107

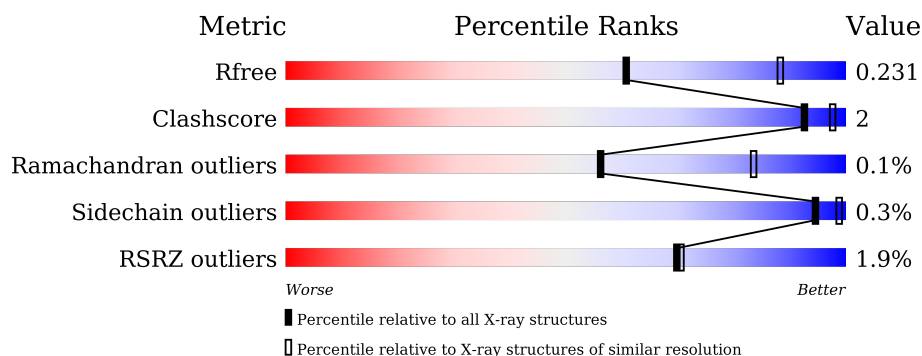
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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. 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	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> % 86% 6% 7% </div> </div>
1	B	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> % 88% 5% 7% </div> </div>
1	C	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 3% 88% . 8% </div> </div>
1	D	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 4% 86% 6% 8% </div> </div>
1	E	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 84% 5% 11% </div> </div>
1	F	338	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; font-size: 0.8em;"> 3% 88% . 9% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14342 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 CRISPR-associated endonuclease Cas1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2379	1504	424	440	11			
1	B	314	Total	C	N	O	S	0	0	0
			2394	1520	417	446	11			
1	C	311	Total	C	N	O	S	0	0	0
			2342	1482	415	434	11			
1	D	312	Total	C	N	O	S	0	0	0
			2330	1486	402	431	11			
1	E	301	Total	C	N	O	S	0	0	0
			2306	1455	408	432	11			
1	F	309	Total	C	N	O	S	0	0	0
			2382	1508	419	444	11			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP Q6D0X0
A	-10	ARG	-	expression tag	UNP Q6D0X0
A	-9	GLY	-	expression tag	UNP Q6D0X0
A	-8	SER	-	expression tag	UNP Q6D0X0
A	-7	HIS	-	expression tag	UNP Q6D0X0
A	-6	HIS	-	expression tag	UNP Q6D0X0
A	-5	HIS	-	expression tag	UNP Q6D0X0
A	-4	HIS	-	expression tag	UNP Q6D0X0
A	-3	HIS	-	expression tag	UNP Q6D0X0
A	-2	HIS	-	expression tag	UNP Q6D0X0
A	-1	GLY	-	expression tag	UNP Q6D0X0
A	0	SER	-	expression tag	UNP Q6D0X0
B	-11	MET	-	initiating methionine	UNP Q6D0X0
B	-10	ARG	-	expression tag	UNP Q6D0X0
B	-9	GLY	-	expression tag	UNP Q6D0X0
B	-8	SER	-	expression tag	UNP Q6D0X0
B	-7	HIS	-	expression tag	UNP Q6D0X0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	HIS	-	expression tag	UNP Q6D0X0
B	-5	HIS	-	expression tag	UNP Q6D0X0
B	-4	HIS	-	expression tag	UNP Q6D0X0
B	-3	HIS	-	expression tag	UNP Q6D0X0
B	-2	HIS	-	expression tag	UNP Q6D0X0
B	-1	GLY	-	expression tag	UNP Q6D0X0
B	0	SER	-	expression tag	UNP Q6D0X0
C	-11	MET	-	initiating methionine	UNP Q6D0X0
C	-10	ARG	-	expression tag	UNP Q6D0X0
C	-9	GLY	-	expression tag	UNP Q6D0X0
C	-8	SER	-	expression tag	UNP Q6D0X0
C	-7	HIS	-	expression tag	UNP Q6D0X0
C	-6	HIS	-	expression tag	UNP Q6D0X0
C	-5	HIS	-	expression tag	UNP Q6D0X0
C	-4	HIS	-	expression tag	UNP Q6D0X0
C	-3	HIS	-	expression tag	UNP Q6D0X0
C	-2	HIS	-	expression tag	UNP Q6D0X0
C	-1	GLY	-	expression tag	UNP Q6D0X0
C	0	SER	-	expression tag	UNP Q6D0X0
D	-11	MET	-	initiating methionine	UNP Q6D0X0
D	-10	ARG	-	expression tag	UNP Q6D0X0
D	-9	GLY	-	expression tag	UNP Q6D0X0
D	-8	SER	-	expression tag	UNP Q6D0X0
D	-7	HIS	-	expression tag	UNP Q6D0X0
D	-6	HIS	-	expression tag	UNP Q6D0X0
D	-5	HIS	-	expression tag	UNP Q6D0X0
D	-4	HIS	-	expression tag	UNP Q6D0X0
D	-3	HIS	-	expression tag	UNP Q6D0X0
D	-2	HIS	-	expression tag	UNP Q6D0X0
D	-1	GLY	-	expression tag	UNP Q6D0X0
D	0	SER	-	expression tag	UNP Q6D0X0
E	-11	MET	-	initiating methionine	UNP Q6D0X0
E	-10	ARG	-	expression tag	UNP Q6D0X0
E	-9	GLY	-	expression tag	UNP Q6D0X0
E	-8	SER	-	expression tag	UNP Q6D0X0
E	-7	HIS	-	expression tag	UNP Q6D0X0
E	-6	HIS	-	expression tag	UNP Q6D0X0
E	-5	HIS	-	expression tag	UNP Q6D0X0
E	-4	HIS	-	expression tag	UNP Q6D0X0
E	-3	HIS	-	expression tag	UNP Q6D0X0
E	-2	HIS	-	expression tag	UNP Q6D0X0
E	-1	GLY	-	expression tag	UNP Q6D0X0

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	SER	-	expression tag	UNP Q6D0X0
F	-11	MET	-	initiating methionine	UNP Q6D0X0
F	-10	ARG	-	expression tag	UNP Q6D0X0
F	-9	GLY	-	expression tag	UNP Q6D0X0
F	-8	SER	-	expression tag	UNP Q6D0X0
F	-7	HIS	-	expression tag	UNP Q6D0X0
F	-6	HIS	-	expression tag	UNP Q6D0X0
F	-5	HIS	-	expression tag	UNP Q6D0X0
F	-4	HIS	-	expression tag	UNP Q6D0X0
F	-3	HIS	-	expression tag	UNP Q6D0X0
F	-2	HIS	-	expression tag	UNP Q6D0X0
F	-1	GLY	-	expression tag	UNP Q6D0X0
F	0	SER	-	expression tag	UNP Q6D0X0

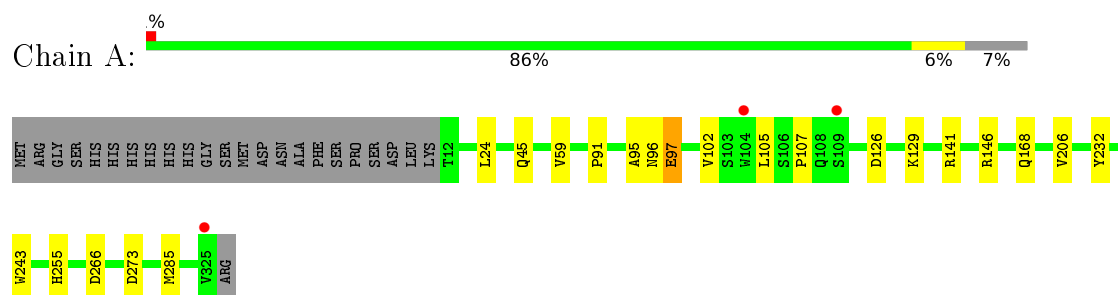
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	44	Total O 44 44	0	0
2	B	33	Total O 33 33	0	0
2	C	30	Total O 30 30	0	0
2	D	22	Total O 22 22	0	0
2	E	55	Total O 55 55	0	0
2	F	25	Total O 25 25	0	0

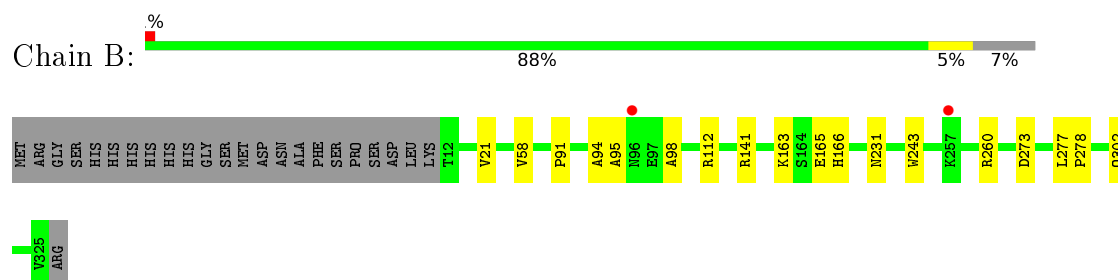
3 Residue-property plots [i](#)

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

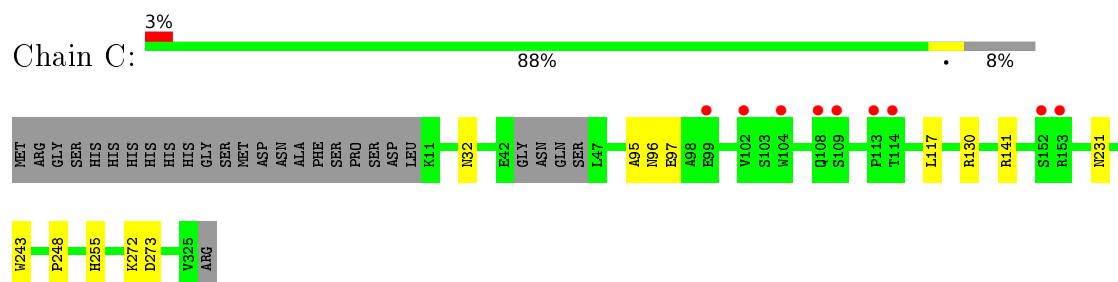
- Molecule 1: CRISPR-associated endonuclease Cas1



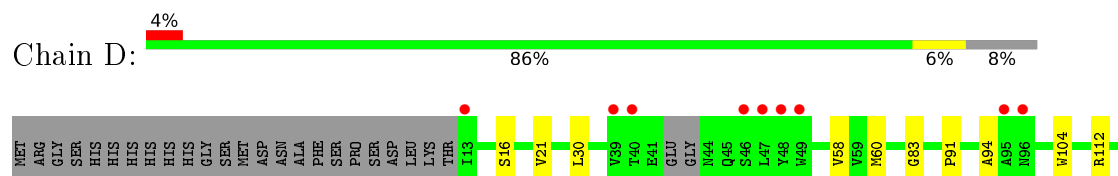
- Molecule 1: CRISPR-associated endonuclease Cas1

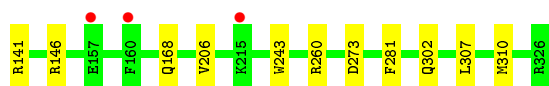


- Molecule 1: CRISPR-associated endonuclease Cas1



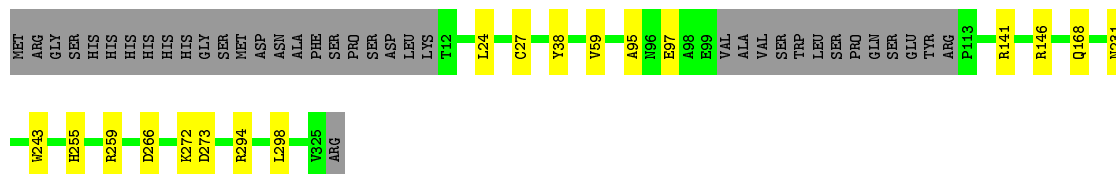
- Molecule 1: CRISPR-associated endonuclease Cas1





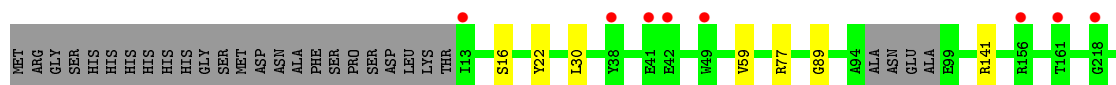
- Molecule 1: CRISPR-associated endonuclease Cas1

Chain E: 84% 5% 11%



- Molecule 1: CRISPR-associated endonuclease Cas1

Chain F: 3% 88% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	261.36 Å 164.06 Å 78.69 Å 90.00° 90.09° 90.00°	Depositor
Resolution (Å)	46.32 – 2.70 46.32 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.8 (46.32-2.70) 90.8 (46.32-2.70)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.69 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.188 , 0.232 0.188 , 0.231	Depositor DCC
R_{free} test set	4008 reflections (4.86%)	DCC
Wilson B-factor (Å ²)	44.3	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.0	EDS
Estimated twinning fraction	0.049 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	1 of 86931 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14342	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.27	0/2426	0.43	0/3296
1	B	0.27	0/2443	0.42	0/3320
1	C	0.26	0/2387	0.43	0/3244
1	D	0.26	0/2378	0.43	0/3238
1	E	0.28	0/2348	0.44	0/3182
1	F	0.27	0/2430	0.43	0/3298
All	All	0.27	0/14412	0.43	0/19578

There are no bond length outliers.

There are no bond angle outliers.

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	2379	0	2269	12	0
1	B	2394	0	2310	11	0
1	C	2342	0	2227	9	0
1	D	2330	0	2199	12	0
1	E	2306	0	2238	10	0
1	F	2382	0	2298	7	0
2	A	44	0	0	0	0
2	B	33	0	0	1	0
2	C	30	0	0	1	0
2	D	22	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	55	0	0	1	0
2	F	25	0	0	0	0
All	All	14342	0	13541	56	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 2.

All (56) 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:E:146:ARG:NH1	1:E:168:GLN:OE1	2.22	0.73
1:D:146:ARG:NH1	1:D:168:GLN:OE1	2.21	0.73
1:E:259:ARG:NH1	2:E:402:HOH:O	2.30	0.64
1:B:163:LYS:HE3	1:B:165:GLU:HB2	1.79	0.64
1:B:141:ARG:NH2	1:B:273:ASP:OD2	2.34	0.61
1:C:32:ASN:ND2	2:C:401:HOH:O	2.34	0.60
1:A:97:GLU:O	1:B:260:ARG:NH2	2.35	0.59
1:A:146:ARG:NH1	1:A:168:GLN:OE1	2.35	0.59
1:F:141:ARG:NH2	1:F:273:ASP:OD2	2.33	0.57
1:D:141:ARG:NH2	1:D:273:ASP:OD2	2.37	0.57
1:D:302:GLN:NE2	2:D:401:HOH:O	2.37	0.57
1:A:255:HIS:ND1	1:A:266:ASP:OD1	2.37	0.55
1:A:96:ASN:O	1:A:243:TRP:NE1	2.38	0.53
1:A:45:GLN:HB3	1:F:77:ARG:HG3	1.90	0.53
1:A:141:ARG:NH2	1:A:273:ASP:OD2	2.44	0.50
1:D:91:PRO:HG2	1:D:94:ALA:HB2	1.93	0.49
1:B:95:ALA:HB3	1:B:98:ALA:HB3	1.93	0.49
1:C:96:ASN:O	1:C:243:TRP:NE1	2.42	0.49
1:D:307:LEU:HA	1:D:310:MET:HE2	1.94	0.49
1:A:206:VAL:HB	1:A:285:MET:HG3	1.94	0.49
1:B:302:GLN:NE2	2:B:401:HOH:O	2.46	0.48
1:A:95:ALA:HB1	1:A:243:TRP:CE2	2.48	0.48
1:B:112:ARG:HD2	1:B:243:TRP:CE2	2.48	0.48
1:B:91:PRO:HG2	1:B:94:ALA:HB2	1.95	0.48
1:D:21:VAL:HG13	1:D:58:VAL:HG13	1.96	0.48
1:D:206:VAL:HG21	1:D:281:PHE:HB3	1.96	0.48
1:F:16:SER:HA	1:F:307:LEU:HD23	1.96	0.48
1:F:307:LEU:HA	1:F:310:MET:HE3	1.97	0.47
1:E:255:HIS:ND1	1:E:266:ASP:OD1	2.39	0.47
1:E:97:GLU:O	1:F:260:ARG:NH2	2.47	0.47
1:C:95:ALA:HB1	1:C:243:TRP:CE2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:ARG:NH2	1:C:273:ASP:OD2	2.48	0.46
1:A:24:LEU:HD13	1:A:59:VAL:HG13	1.97	0.46
1:A:105:LEU:O	1:A:107:PRO:HD3	2.16	0.45
1:E:95:ALA:HB1	1:E:243:TRP:CE2	2.51	0.45
1:B:165:GLU:OE1	1:B:165:GLU:N	2.48	0.45
1:E:294:ARG:NH1	1:E:298:LEU:HD11	2.32	0.44
1:B:21:VAL:HG13	1:B:58:VAL:HG13	1.99	0.44
1:E:231:ASN:CG	1:E:272:LYS:HD2	2.37	0.44
1:E:27:CYS:SG	1:E:38:TYR:HB3	2.58	0.44
1:F:89:GLY:HA3	1:F:228:ASP:HB3	2.00	0.44
1:D:16:SER:HA	1:D:307:LEU:HD23	2.00	0.44
1:C:96:ASN:ND2	1:D:104:TRP:O	2.43	0.43
1:D:60:MET:HE1	1:D:83:GLY:HA3	2.00	0.43
1:D:112:ARG:HD2	1:D:243:TRP:CE2	2.53	0.43
1:E:141:ARG:NH2	1:E:273:ASP:OD2	2.51	0.43
1:A:91:PRO:HG3	1:A:232:TYR:CZ	2.54	0.43
1:C:231:ASN:CG	1:C:272:LYS:HD2	2.39	0.42
1:C:117:LEU:HD13	1:C:248:PRO:HD3	2.02	0.42
1:C:97:GLU:O	1:D:260:ARG:NH2	2.51	0.42
1:A:126:ASP:HB3	1:A:129:LYS:HB2	2.02	0.42
1:F:22:TYR:HB2	1:F:59:VAL:HG22	2.02	0.42
1:B:163:LYS:HB3	1:B:166:HIS:CD2	2.55	0.42
1:C:130:ARG:NH2	1:C:255:HIS:O	2.54	0.41
1:B:277:LEU:HB3	1:B:278:PRO:HD3	2.03	0.41
1:E:24:LEU:HD13	1:E:59:VAL:HG13	2.03	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	312/338 (92%)	303 (97%)	8 (3%)	1 (0%)	46 75

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	312/338 (92%)	306 (98%)	6 (2%)	0	100	100
1	C	307/338 (91%)	303 (99%)	4 (1%)	0	100	100
1	D	308/338 (91%)	300 (97%)	8 (3%)	0	100	100
1	E	297/338 (88%)	294 (99%)	3 (1%)	0	100	100
1	F	305/338 (90%)	302 (99%)	3 (1%)	0	100	100
All	All	1841/2028 (91%)	1808 (98%)	32 (2%)	1 (0%)	56	83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	102	VAL

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	232/275 (84%)	231 (100%)	1 (0%)	93	98
1	B	240/275 (87%)	239 (100%)	1 (0%)	93	98
1	C	227/275 (82%)	227 (100%)	0	100	100
1	D	224/275 (82%)	223 (100%)	1 (0%)	93	98
1	E	233/275 (85%)	233 (100%)	0	100	100
1	F	241/275 (88%)	240 (100%)	1 (0%)	93	98
All	All	1397/1650 (85%)	1393 (100%)	4 (0%)	94	99

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	GLU
1	B	231	ASN
1	D	30	LEU
1	F	30	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	249	HIS

5.3.3 RNA [i](#)

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	314/338 (92%)	-0.21	3 (0%) 84 85	29, 46, 80, 95	0
1	B	314/338 (92%)	-0.15	2 (0%) 90 91	29, 50, 83, 111	0
1	C	311/338 (92%)	-0.01	9 (2%) 55 55	34, 52, 83, 108	0
1	D	312/338 (92%)	0.13	12 (3%) 44 44	33, 61, 90, 116	0
1	E	301/338 (89%)	-0.35	0 100 100	26, 40, 65, 85	0
1	F	309/338 (91%)	-0.01	9 (2%) 55 55	29, 48, 89, 107	0
All	All	1861/2028 (91%)	-0.10	35 (1%) 70 70	26, 49, 84, 116	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	102	VAL	4.5
1	A	104	TRP	3.9
1	F	42	GLU	3.5
1	F	156	ARG	3.5
1	D	96	ASN	3.4
1	F	13	ILE	3.3
1	D	49	TRP	3.2
1	C	109	SER	3.2
1	D	95	ALA	3.2
1	A	109	SER	3.0
1	C	108	GLN	3.0
1	A	325	VAL	3.0
1	F	49	TRP	2.9
1	C	99	GLU	2.8
1	D	39	VAL	2.7
1	D	157	GLU	2.7
1	C	152	SER	2.6
1	D	215	LYS	2.6
1	D	47	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	96	ASN	2.5
1	C	104	TRP	2.5
1	D	48	TYR	2.5
1	F	41	GLU	2.5
1	F	38	TYR	2.4
1	D	46	SER	2.4
1	B	257	LYS	2.4
1	D	40	THR	2.3
1	F	258	THR	2.2
1	F	161	THR	2.2
1	D	160	PHE	2.2
1	F	218	GLY	2.2
1	C	113	PRO	2.2
1	C	114	THR	2.1
1	D	13	ILE	2.1
1	C	153	ARG	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.