Full wwPDB X-ray Structure Validation Report ⓘ

Sep 25, 2023 – 11:04 AM EDT

PDB ID	:	5VVK
Title	:	Cas1-Cas2 bound to full-site mimic
Authors	:	Wright, A.V.; Knott, G.J.; Doxzen, K.D.; Doudna, J.A.
Deposited on	:	2017-05-19
Resolution	:	2.90 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

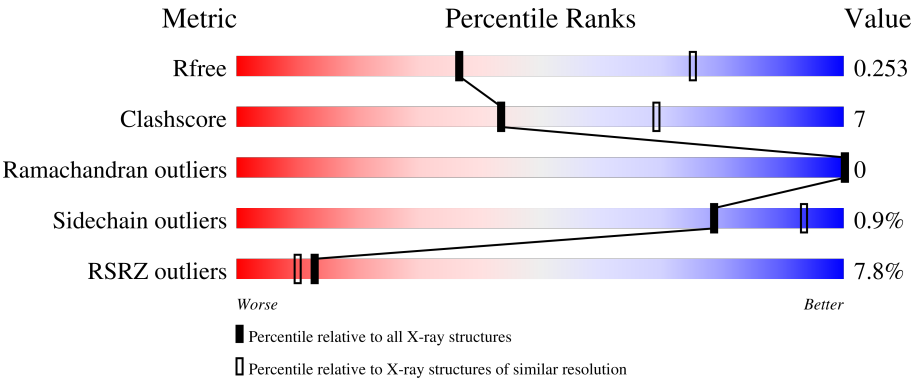
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	308	11% 76% 10% 14%
1	B	308	11% 72% 13% 15%
1	C	308	5% 72% 14% 13%
1	D	308	5% 76% 13% 12%
2	E	94	6% 83% 17%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	94	4% 81% 18%
3	G	11	64% 36%
4	H	11	36% 64%
5	J	58	21% 53% 26%
6	K	58	34% 40% 26%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 11896 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	266	Total	C	N	O	S	0	0	0
			2045	1306	363	369	7			
1	B	263	Total	C	N	O	S	0	0	0
			2017	1290	361	359	7			
1	C	267	Total	C	N	O	S	0	0	0
			2052	1311	364	370	7			
1	D	272	Total	C	N	O	S	0	0	0
			2096	1340	372	377	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q46896
A	-1	PHE	-	expression tag	UNP Q46896
A	0	THR	-	expression tag	UNP Q46896
B	-2	SER	-	expression tag	UNP Q46896
B	-1	PHE	-	expression tag	UNP Q46896
B	0	THR	-	expression tag	UNP Q46896
C	-2	SER	-	expression tag	UNP Q46896
C	-1	PHE	-	expression tag	UNP Q46896
C	0	THR	-	expression tag	UNP Q46896
D	-2	SER	-	expression tag	UNP Q46896
D	-1	PHE	-	expression tag	UNP Q46896
D	0	THR	-	expression tag	UNP Q46896

- Molecule 2 is a protein called CRISPR-associated endoribonuclease Cas2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	94	Total	C	N	O	S	0	0	0
			739	475	128	132	4			
2	F	94	Total	C	N	O	S	0	0	0
			739	475	128	132	4			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*CP*CP*CP*CP*AP*GP*TP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	11	Total	C	N	O	P	0	0	0
			220	105	42	63	10			

- Molecule 4 is a DNA chain called DNA (5'-D(*GP*AP*CP*CP*AP*CP*CP*AP*GP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	11	Total	C	N	O	P	0	0	0
			222	106	44	62	10			

- Molecule 5 is a DNA chain called DNA (58-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	J	43	Total	C	N	O	P	0	0	0
			883	421	164	256	42			

- Molecule 6 is a DNA chain called DNA (58-MER).

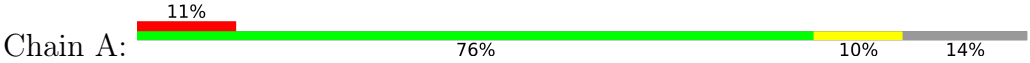
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	K	43	Total	C	N	O	P	0	0	0
			883	419	166	256	42			

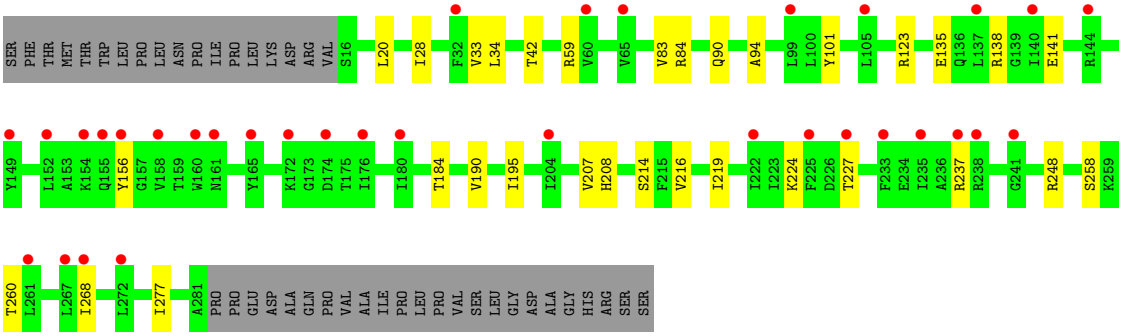
3 Residue-property plots (i)

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

• Molecule 1: CRISPR-associated endonuclease Cas1

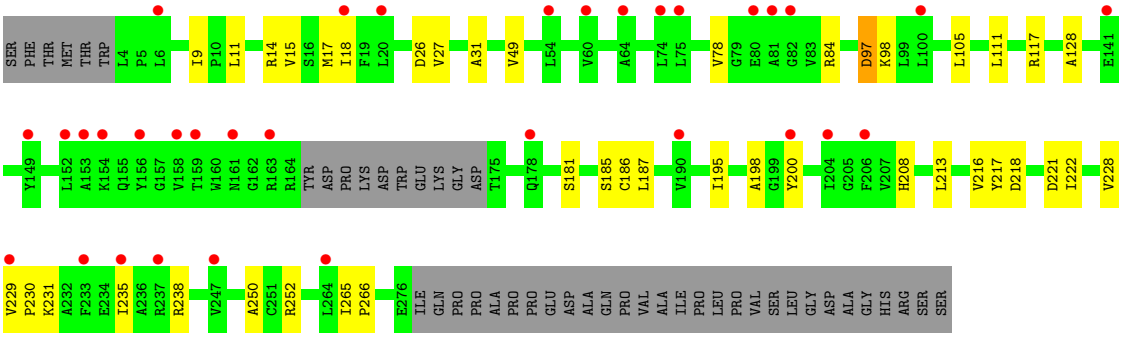
Chain A:





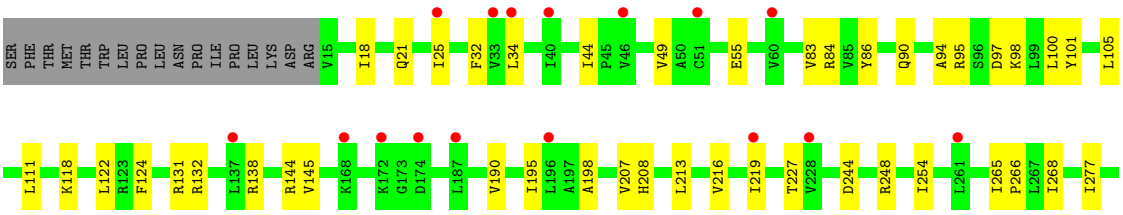
• Molecule 1: CRISPR-associated endonuclease Cas1

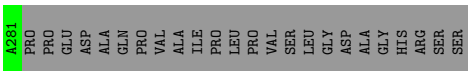
Chain B:



• Molecule 1: CRISPR-associated endonuclease Cas1

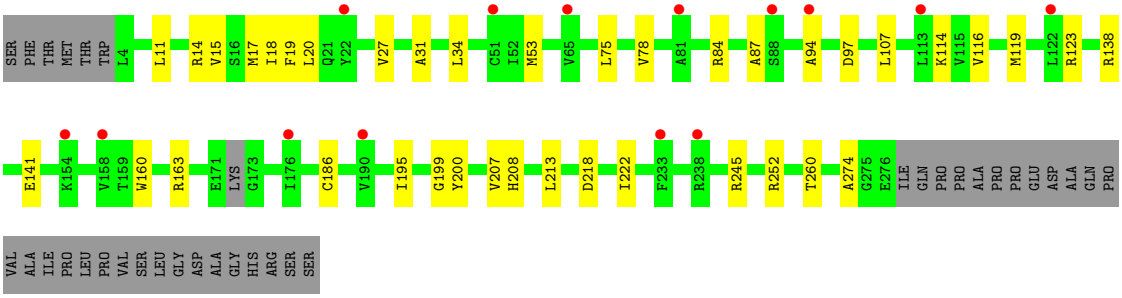
Chain C:





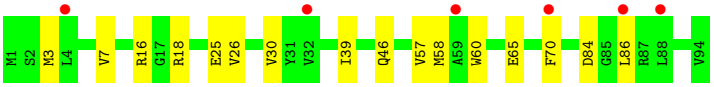
• Molecule 1: CRISPR-associated endonuclease Cas1

Chain D:



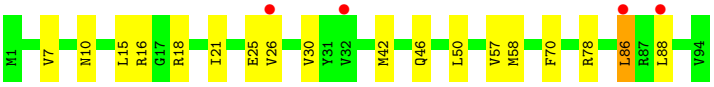
• Molecule 2: CRISPR-associated endoribonuclease Cas2

Chain E:



• Molecule 2: CRISPR-associated endoribonuclease Cas2

Chain F:

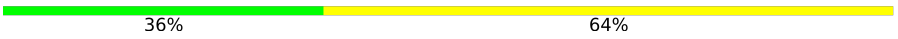


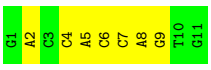
• Molecule 3: DNA (5'-D(*GP*CP*CP*CP*CP*AP*GP*TP*AP*GP*C)-3')

Chain G:

• Molecule 4: DNA (5'-D(*GP*AP*CP*CP*AP*CP*CP*AP*GP*TP*G)-3')

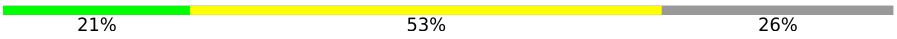
Chain H:





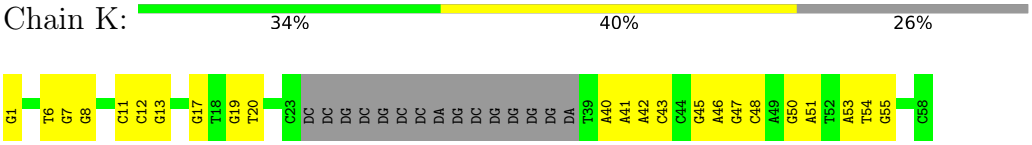
• Molecule 5: DNA (58-MER)

Chain J:



• Molecule 6: DNA (58-MER)

Chain K:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	74.90Å 197.61Å 95.34Å 90.00° 112.70° 90.00°	Depositor
Resolution (Å)	98.81 – 2.90 98.81 – 2.76	Depositor EDS
% Data completeness (in resolution range)	99.8 (98.81-2.90) 98.1 (98.81-2.76)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.14 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.11.1	Depositor
R, R_{free}	0.217 , 0.254 0.216 , 0.253	Depositor DCC
R_{free} test set	3349 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å^2)	72.3	Xtriage
Anisotropy	0.708	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 67.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	11896	wwPDB-VP
Average B, all atoms (Å^2)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.24	0/2085	0.39	0/2827
1	B	0.24	0/2053	0.39	0/2782
1	C	0.25	0/2092	0.41	0/2837
1	D	0.25	0/2136	0.41	0/2896
2	E	0.26	0/753	0.48	0/1024
2	F	0.25	0/753	0.47	0/1024
3	G	0.55	0/246	0.81	0/377
4	H	0.55	0/249	0.79	0/382
5	J	0.53	0/990	0.91	0/1525
6	K	0.57	0/990	0.90	0/1525
All	All	0.33	0/12347	0.55	0/17199

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	2045	0	2097	17	0
1	B	2017	0	2095	27	0
1	C	2052	0	2106	26	0
1	D	2096	0	2155	25	0
2	E	739	0	756	15	0
2	F	739	0	756	16	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	220	0	124	4	0
4	H	222	0	124	4	0
5	J	883	0	487	25	0
6	K	883	0	485	18	0
All	All	11896	0	11185	155	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 7.

All (155) 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:D:18:ILE:HD11	2:E:86:LEU:HD23	1.64	0.80
5:J:39:DG:H2"	5:J:40:DA:C8	2.17	0.79
5:J:47:DT:H2"	5:J:48:DA:C8	2.23	0.74
1:C:145:VAL:HG21	5:J:58:DA:H5'	1.71	0.72
1:A:59:ARG:NH1	4:H:2:DA:OP1	2.23	0.72
1:D:123:ARG:NH1	1:D:141:GLU:OE1	2.23	0.71
2:E:18:ARG:NH1	2:E:46:GLN:OE1	2.24	0.71
1:B:217:TYR:O	1:B:221:ASP:HB2	1.92	0.69
1:C:18:ILE:HG22	1:C:49:VAL:HG21	1.78	0.65
1:D:138:ARG:NH1	1:D:207:VAL:O	2.29	0.63
5:J:46:DC:H2'	5:J:47:DT:H71	1.82	0.62
1:B:97:ASP:OD1	1:B:97:ASP:N	2.32	0.61
1:D:114:LYS:HB3	1:D:274:ALA:HB1	1.83	0.61
2:F:16:ARG:NH2	2:F:25:GLU:OE2	2.28	0.61
6:K:54:DT:H2"	6:K:55:DG:H8	1.67	0.60
2:E:16:ARG:NH2	2:E:25:GLU:OE2	2.24	0.60
1:D:27:VAL:HA	1:D:31:ALA:O	2.01	0.60
1:A:59:ARG:NH2	1:B:26:ASP:OD1	2.35	0.60
1:C:90:GLN:HB3	1:C:94:ALA:HB2	1.85	0.59
1:C:208:HIS:NE2	6:K:17:DG:OP2	2.35	0.59
1:D:11:LEU:HD12	1:D:14:ARG:HD2	1.85	0.58
5:J:17:DG:H2'	5:J:18:DG:C8	2.38	0.58
1:C:248:ARG:NH2	6:K:11:DC:OP1	2.36	0.57
1:C:84:ARG:NH1	6:K:13:DG:OP1	2.37	0.57
6:K:45:DG:H2'	6:K:46:DA:C8	2.39	0.57
1:C:25:ILE:HG13	1:C:34:LEU:HD23	1.87	0.57
1:A:156:TYR:HD1	1:A:237:ARG:HD2	1.71	0.56
1:A:184:THR:HG22	1:A:224:LYS:HD2	1.86	0.56
2:E:7:VAL:HG21	2:F:30:VAL:HG11	1.88	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:K:46:DA:H2'	6:K:47:DG:C8	2.41	0.56
2:F:18:ARG:NH1	2:F:46:GLN:OE1	2.39	0.55
5:J:10:DT:H2'	5:J:11:DC:C6	2.41	0.55
1:B:198:ALA:HB2	1:B:265:ILE:HD12	1.87	0.55
1:D:94:ALA:HA	1:D:199:GLY:HA2	1.87	0.55
5:J:9:DG:H2'	5:J:10:DT:H71	1.89	0.54
5:J:38:DG:H2"	5:J:39:DG:C8	2.42	0.54
1:A:83:VAL:HG12	1:A:84:ARG:HG2	1.89	0.54
1:B:181:SER:O	1:B:185:SER:HB2	2.08	0.53
5:J:4:DT:H2"	5:J:5:DG:C8	2.44	0.53
6:K:19:DG:H4'	6:K:20:DT:OP1	2.09	0.53
1:A:20:LEU:HD22	1:A:34:LEU:HD22	1.91	0.52
1:A:138:ARG:NH1	1:A:207:VAL:O	2.41	0.52
5:J:55:DT:H2"	5:J:56:DA:C8	2.45	0.52
1:B:9:ILE:HG12	2:E:39:ILE:HD12	1.92	0.52
1:B:17:MET:O	1:B:252:ARG:NH1	2.43	0.52
5:J:4:DT:H2"	5:J:5:DG:H8	1.76	0.51
1:A:123:ARG:NH2	1:A:141:GLU:OE2	2.43	0.50
1:C:21:GLN:HB3	1:C:55:GLU:HB2	1.92	0.50
6:K:54:DT:H2"	6:K:55:DG:C8	2.45	0.50
1:C:83:VAL:HG12	1:C:84:ARG:HG2	1.91	0.50
1:C:131:ARG:O	1:C:132:ARG:NH1	2.40	0.50
3:G:5:DC:H2"	3:G:6:DA:C8	2.46	0.50
1:A:248:ARG:NH1	5:J:11:DC:OP1	2.41	0.50
1:B:187:LEU:HD22	1:B:228:VAL:HG21	1.94	0.50
2:F:78:ARG:NH1	5:J:3:DC:OP1	2.41	0.50
1:B:15:VAL:HG11	2:F:86:LEU:HD13	1.93	0.49
1:B:231:LYS:HG2	1:B:250:ALA:HB1	1.93	0.49
5:J:1:DC:H2"	5:J:2:DA:C8	2.47	0.49
1:B:195:ILE:HD12	1:B:216:VAL:HG22	1.93	0.49
2:F:86:LEU:HD23	2:F:88:LEU:HG	1.95	0.49
2:E:30:VAL:HG11	2:F:7:VAL:HG21	1.94	0.49
6:K:47:DG:H2'	6:K:48:DC:C6	2.48	0.49
1:B:9:ILE:HB	1:B:14:ARG:HH21	1.78	0.48
1:B:11:LEU:HD12	1:B:14:ARG:HD2	1.95	0.48
6:K:6:DT:H2"	6:K:7:DG:C8	2.48	0.48
1:D:20:LEU:HD22	1:D:34:LEU:HD22	1.94	0.48
3:G:8:DT:H2"	3:G:9:DA:H8	1.79	0.48
1:C:118:LYS:HG3	1:C:122:LEU:HD12	1.95	0.48
1:B:208:HIS:ND1	1:B:218:ASP:OD1	2.41	0.48
1:D:195:ILE:HG23	1:D:200:TYR:HB2	1.96	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:J:40:DA:H1'	5:J:41:DA:C8	2.47	0.48
1:B:218:ASP:O	1:B:222:ILE:HG13	2.14	0.47
2:F:46:GLN:O	2:F:50:LEU:HB2	2.14	0.47
1:C:124:PHE:CE1	1:C:144:ARG:HD2	2.49	0.47
1:B:18:ILE:HG12	1:B:49:VAL:HG21	1.96	0.47
1:C:198:ALA:HB2	1:C:265:ILE:HG23	1.97	0.47
1:D:116:VAL:HG13	1:D:207:VAL:HG22	1.96	0.47
5:J:42:DC:H4'	5:J:43:DA:OP1	2.14	0.47
1:D:97:ASP:OD1	1:D:97:ASP:N	2.48	0.46
1:D:245:ARG:HG3	2:E:84:ASP:HA	1.97	0.46
2:E:26:VAL:HG21	2:F:57:VAL:HG21	1.97	0.46
5:J:45:DT:H5'	5:J:45:DT:H6	1.81	0.46
5:J:55:DT:H2"	5:J:56:DA:N7	2.30	0.46
5:J:3:DC:C6	5:J:4:DT:H72	2.50	0.46
1:C:138:ARG:NH2	1:C:207:VAL:O	2.48	0.46
1:D:84:ARG:CZ	1:D:213:LEU:HD11	2.46	0.45
1:B:17:MET:HE1	1:B:186:CYS:HB3	1.97	0.45
1:C:101:TYR:HE1	1:C:277:ILE:HG22	1.80	0.45
1:D:252:ARG:NH2	2:E:84:ASP:O	2.40	0.45
2:E:86:LEU:HA	2:E:86:LEU:HD12	1.70	0.45
1:C:195:ILE:HD12	1:C:216:VAL:HG22	1.98	0.45
1:D:75:LEU:O	1:D:87:ALA:HA	2.16	0.45
1:A:195:ILE:HD12	1:A:216:VAL:HG22	1.98	0.45
6:K:7:DG:H2"	6:K:8:DG:H8	1.82	0.45
4:H:4:DC:H2"	4:H:5:DA:C8	2.52	0.44
2:E:57:VAL:HG21	2:F:26:VAL:HG21	1.98	0.44
1:C:32:PHE:HB2	1:C:44:ILE:HB	1.99	0.44
1:C:86:TYR:OH	6:K:12:DC:OP2	2.20	0.44
6:K:40:DA:H2"	6:K:41:DA:C8	2.52	0.44
6:K:42:DA:H2"	6:K:43:DC:OP2	2.17	0.44
1:C:244:ASP:N	1:C:244:ASP:OD1	2.51	0.44
1:D:11:LEU:HA	1:D:14:ARG:HD2	2.00	0.44
1:C:100:LEU:HB3	1:D:107:LEU:HD21	1.99	0.44
1:B:117:ARG:HG3	1:B:128:ALA:HB3	1.99	0.44
1:D:160:TRP:CZ2	1:D:163:ARG:HB2	2.53	0.44
1:A:28:ILE:HG13	1:A:33:VAL:HG11	2.00	0.44
1:A:90:GLN:HG2	1:A:94:ALA:HB2	2.00	0.44
1:A:219:ILE:HG12	1:A:268:ILE:HG12	2.00	0.43
1:B:27:VAL:HA	1:B:31:ALA:O	2.18	0.43
1:D:19:PHE:CZ	1:D:53:MET:HG3	2.53	0.43
1:D:208:HIS:ND1	1:D:218:ASP:OD1	2.42	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:10:ASN:HB2	6:K:1:DG:H5"	2.00	0.43
1:C:265:ILE:HB	1:C:266:PRO:HD3	1.99	0.43
2:F:18:ARG:HD3	2:F:18:ARG:HA	1.86	0.43
1:B:78:VAL:HG12	1:B:84:ARG:HA	2.00	0.43
1:D:119:MET:HG2	1:D:222:ILE:HD11	2.00	0.43
2:F:42:MET:O	2:F:46:GLN:HG2	2.19	0.43
5:J:43:DA:C2	6:K:19:DG:C2	3.07	0.43
1:C:213:LEU:HD22	1:C:216:VAL:HG21	2.01	0.43
1:D:84:ARG:HH21	1:D:87:ALA:HB3	1.84	0.43
1:B:105:LEU:HD23	1:B:111:LEU:HD13	2.01	0.43
1:C:219:ILE:HG12	1:C:268:ILE:HG12	2.01	0.43
1:C:227:THR:HG22	1:C:254:ILE:HD13	2.01	0.43
1:D:15:VAL:HG13	2:E:65:GLU:OE1	2.19	0.42
6:K:50:DG:H2"	6:K:51:DA:C8	2.53	0.42
1:D:17:MET:HE1	1:D:186:CYS:HB3	2.01	0.42
5:J:13:DC:H4'	5:J:14:DC:C5	2.54	0.42
1:B:265:ILE:HB	1:B:266:PRO:HD3	2.02	0.42
1:B:98:LYS:HB3	1:B:200:TYR:CE1	2.54	0.42
2:E:3:MET:O	2:E:60:TRP:HA	2.20	0.42
1:C:95:ARG:NH1	1:C:98:LYS:HE3	2.35	0.42
4:H:8:DA:H2"	4:H:9:DG:C8	2.54	0.42
1:A:208:HIS:O	1:A:214:SER:HB3	2.20	0.42
5:J:7:DT:H2"	5:J:8:DG:C8	2.54	0.42
1:A:34:LEU:HB3	1:A:42:THR:HB	2.01	0.42
3:G:5:DC:H2"	3:G:6:DA:H8	1.84	0.42
1:D:218:ASP:O	1:D:222:ILE:HG13	2.20	0.41
2:E:58:MET:O	2:E:70:PHE:HA	2.20	0.41
2:F:58:MET:O	2:F:70:PHE:HA	2.20	0.41
5:J:50:DG:N2	5:J:52:DT:H1'	2.36	0.41
5:J:21:DT:H2"	5:J:22:DA:C8	2.55	0.41
2:F:18:ARG:HD3	2:F:21:ILE:HD12	2.02	0.41
4:H:6:DC:H2"	4:H:7:DC:C6	2.55	0.41
1:A:101:TYR:HE1	1:A:277:ILE:HG22	1.86	0.41
1:C:105:LEU:HD23	1:C:111:LEU:HD13	2.02	0.41
1:B:229:VAL:N	1:B:230:PRO:HD2	2.35	0.41
1:B:235:ILE:HD12	1:B:238:ARG:HH21	1.86	0.41
5:J:40:DA:H4'	5:J:41:DA:OP1	2.21	0.41
1:B:217:TYR:O	1:B:221:ASP:CB	2.67	0.41
2:E:7:VAL:HB	2:E:57:VAL:HG13	2.03	0.40
2:F:15:LEU:HB2	2:F:50:LEU:HB3	2.03	0.40
6:K:53:DA:H2"	6:K:54:DT:O4'	2.22	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:SER:HB2	1:A:260:THR:HG23	2.03	0.40
1:B:213:LEU:HB2	1:B:217:TYR:CZ	2.57	0.40
3:G:8:DT:H2"	3:G:9:DA:C8	2.56	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	264/308 (86%)	262 (99%)	2 (1%)	0	100	100
1	B	259/308 (84%)	253 (98%)	6 (2%)	0	100	100
1	C	265/308 (86%)	260 (98%)	5 (2%)	0	100	100
1	D	268/308 (87%)	262 (98%)	6 (2%)	0	100	100
2	E	92/94 (98%)	89 (97%)	3 (3%)	0	100	100
2	F	92/94 (98%)	88 (96%)	4 (4%)	0	100	100
All	All	1240/1420 (87%)	1214 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	211/248 (85%)	208 (99%)	3 (1%)	67 89
1	B	210/248 (85%)	209 (100%)	1 (0%)	88 96
1	C	212/248 (86%)	210 (99%)	2 (1%)	78 93
1	D	218/248 (88%)	216 (99%)	2 (1%)	78 93
2	E	79/79 (100%)	79 (100%)	0	100 100
2	F	79/79 (100%)	78 (99%)	1 (1%)	69 90
All	All	1009/1150 (88%)	1000 (99%)	9 (1%)	78 93

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	GLU
1	A	190	VAL
1	A	227	THR
1	B	97	ASP
1	C	97	ASP
1	C	190	VAL
1	D	78	VAL
1	D	260	THR
2	F	86	LEU

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

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 monosaccharides 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, 95^{th} 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	266/308 (86%)	0.81	34 (12%) 3 2	59, 99, 139, 158	0
1	B	263/308 (85%)	0.92	33 (12%) 3 3	55, 105, 164, 171	0
1	C	267/308 (86%)	0.69	16 (5%) 21 18	57, 85, 116, 157	0
1	D	272/308 (88%)	0.67	14 (5%) 28 24	47, 85, 125, 149	0
2	E	94/94 (100%)	0.72	6 (6%) 19 15	48, 70, 96, 108	0
2	F	94/94 (100%)	0.53	4 (4%) 35 31	52, 68, 101, 116	0
3	G	11/11 (100%)	-0.03	0 100 100	63, 88, 177, 189	0
4	H	11/11 (100%)	0.06	0 100 100	68, 93, 199, 202	0
5	J	43/58 (74%)	-0.37	0 100 100	74, 129, 191, 208	0
6	K	43/58 (74%)	-0.25	0 100 100	67, 145, 198, 210	0
All	All	1364/1558 (87%)	0.67	107 (7%) 13 10	47, 90, 158, 210	0

All (107) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	152	LEU	10.3
1	A	158	VAL	5.0
1	A	272	LEU	4.9
1	A	156	TYR	4.7
1	A	155	GLN	4.6
1	B	153	ALA	4.4
1	A	152	LEU	4.4
1	D	238	ARG	4.2
1	B	233	PHE	4.1
1	B	159	THR	4.0
1	B	149	TYR	4.0
1	A	99	LEU	3.6
1	B	229	VAL	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	6	LEU	3.5
1	C	46	VAL	3.4
1	A	225	PHE	3.4
1	A	176	ILE	3.3
1	A	180	ILE	3.3
1	B	237	ARG	3.3
1	B	154	LYS	3.3
1	B	141	GLU	3.2
1	B	82	GLY	3.1
1	C	137	LEU	3.0
1	C	261	LEU	3.0
1	B	161	ASN	3.0
1	B	156	TYR	2.9
1	A	227	THR	2.9
1	B	80	GLU	2.8
1	B	158	VAL	2.8
1	A	238	ARG	2.7
1	A	235	ILE	2.7
1	A	137	LEU	2.7
1	A	233	PHE	2.7
1	D	94	ALA	2.7
1	D	122	LEU	2.6
1	B	18	ILE	2.6
2	E	4	LEU	2.6
1	B	100	LEU	2.6
1	B	204	ILE	2.6
1	B	264	LEU	2.6
1	D	158	VAL	2.5
1	C	196	LEU	2.5
1	C	60	VAL	2.5
1	B	54	LEU	2.5
1	A	204	ILE	2.5
1	D	190	VAL	2.4
1	B	20	LEU	2.4
1	B	200	TYR	2.4
2	F	88	LEU	2.4
1	D	233	PHE	2.4
1	D	65	VAL	2.4
1	D	88	SER	2.4
1	C	34	LEU	2.4
2	E	88	LEU	2.4
1	A	32	PHE	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	22	TYR	2.4
1	A	261	LEU	2.4
1	B	81	ALA	2.3
1	C	168	LYS	2.3
1	B	75	LEU	2.3
1	A	222	ILE	2.3
1	D	113	LEU	2.3
1	A	237	ARG	2.3
1	A	140	ILE	2.3
1	B	235	ILE	2.3
1	C	219	ILE	2.3
1	C	33	VAL	2.3
1	A	154	LYS	2.3
1	A	172	LYS	2.3
1	B	247	VAL	2.2
2	E	86	LEU	2.2
1	A	268	ILE	2.2
1	A	149	TYR	2.2
2	F	26	VAL	2.2
2	E	32	VAL	2.2
1	A	144	ARG	2.2
1	B	64	ALA	2.2
1	B	60	VAL	2.2
1	A	267	LEU	2.2
2	F	32	VAL	2.1
1	C	40	ILE	2.1
1	C	51	CYS	2.1
1	D	51	CYS	2.1
1	A	105	LEU	2.1
1	B	163	ARG	2.1
1	A	160	TRP	2.1
1	A	165	TYR	2.1
1	A	60	VAL	2.1
1	C	228	VAL	2.1
1	D	154	LYS	2.1
2	E	70	PHE	2.1
1	B	178	GLN	2.0
1	A	174	ASP	2.0
1	A	161	ASN	2.0
1	D	176	ILE	2.0
1	A	65	VAL	2.0
1	B	206	PHE	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	172	LYS	2.0
1	C	174	ASP	2.0
1	B	190	VAL	2.0
1	B	74	LEU	2.0
1	C	187	LEU	2.0
2	F	86	LEU	2.0
1	A	241	GLY	2.0
1	C	25	ILE	2.0
1	D	81	ALA	2.0
2	E	59	ALA	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 monosaccharides 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.