Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:20 am BST

PDB ID	:	5H58
Title	:	Structural and dynamics studies of the TetR family protein, CprB from Streptomyces coelicolor in complex with its biological operator sequence
Authors	:	Bhukya, H.; Jana, A.K.; Sengupta, N.; Anand, R.
Deposited on	:	2016-11-04
Resolution	:	3.99 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 following versions of software and data (see references ⓘ) were used in the production of this report:

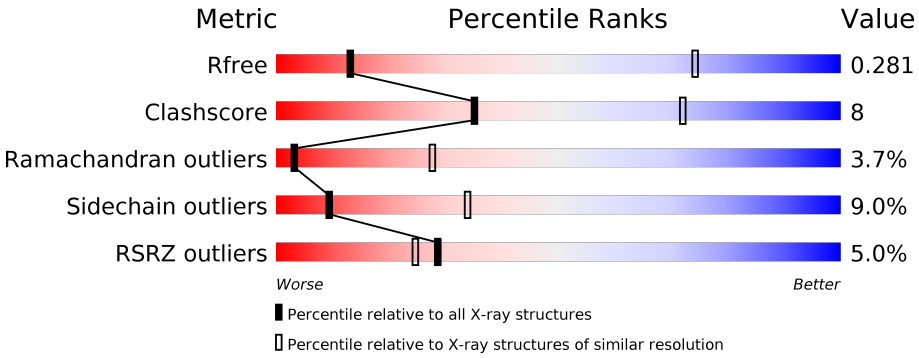
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 3.99 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	1087 (4.30-3.70)
Clashscore	141614	1148 (4.30-3.70)
Ramachandran outliers	138981	1108 (4.30-3.70)
Sidechain outliers	138945	1099 (4.30-3.70)
RSRZ outliers	127900	1028 (4.34-3.66)

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 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	215	2% 73% 16% • 10%
1	B	215	5% 73% 18% • 5%
1	C	215	5% 70% 22% • 7%
1	D	215	8% 62% 25% • 10%
2	E	27	30% 44% 26%
3	F	27	4% 48% 33% 19%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6885 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 CprB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	194	Total	C	N	O	S	0	0	0
			1480	930	276	269	5			
1	B	204	Total	C	N	O	S	0	0	0
			1518	950	277	285	6			
1	C	200	Total	C	N	O	S	0	0	0
			1535	961	295	274	5			
1	D	193	Total	C	N	O	S	0	0	0
			1497	939	279	273	6			

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*GP*GP*C*AP*GP*GP*CP*GP*GP*CP*AP*CP*GP*GP*TP*CP*TP*GP*TP*TP*GP*AP*GP*TP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	20	Total	C	N	O	P	0	0	0
			414	196	80	119	19			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*AP*A*CP*TP*CP*AP*AP*CP*AP*GP*AP*CP*CP*GP*TP*GP*CP*CP*GP*CP*CP*TP*GP*CP*CP*T)-3').

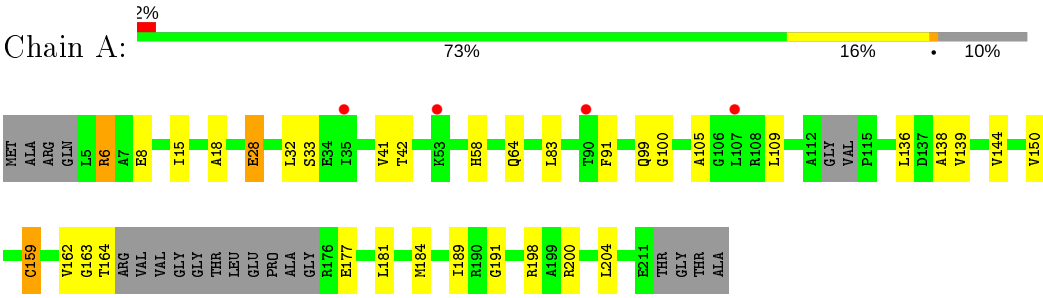
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	22	Total	C	N	O	P	0	0	0
			441	210	81	129	21			

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.

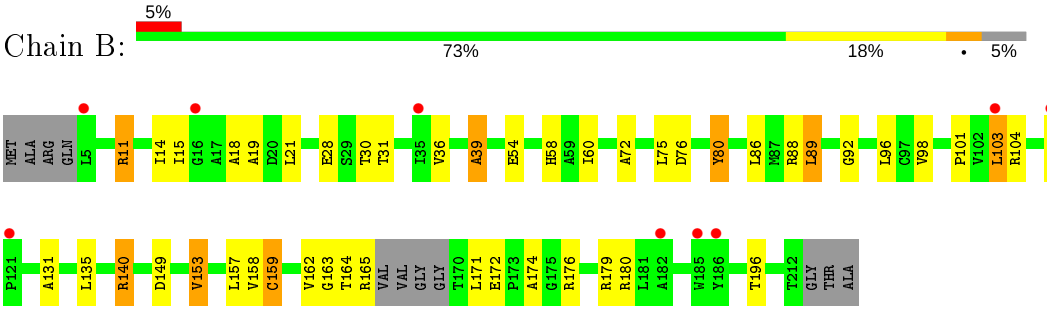
- Molecule 1: CprB

Chain A:



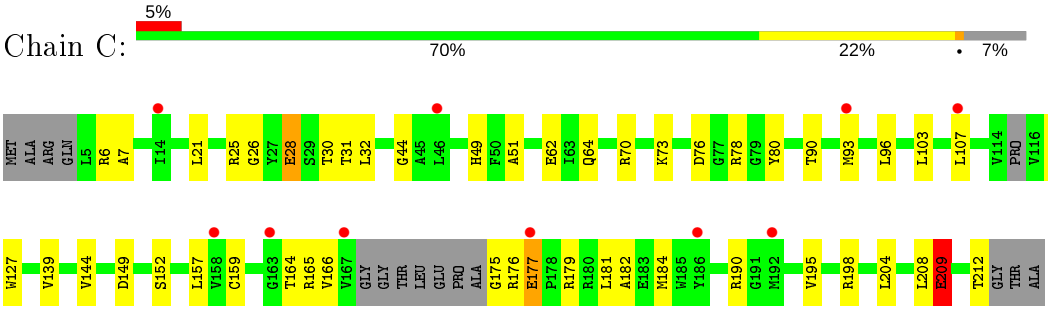
- Molecule 1: CprB

Chain B:



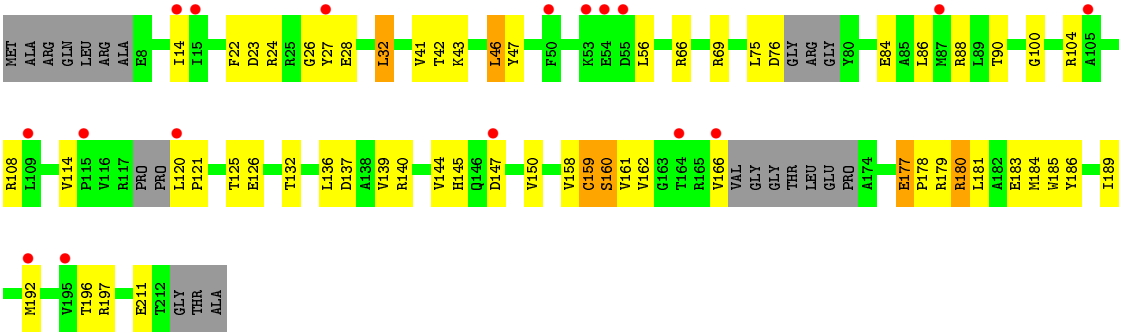
- Molecule 1: CprB

Chain C:



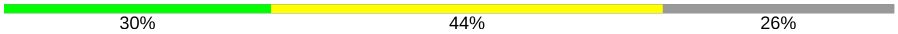
- Molecule 1: CprB

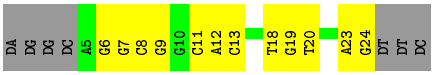
Chain D:



• Molecule 2: DNA (5'-D(*AP*GP*GP*C*AP*GP*GP*CP*GP*GP*CP*AP*CP*GP*GP*TP*CP*TP*GP*TP*TP*GP*AP*GP*TP*TP*C)-3')

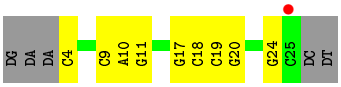
Chain E:





• Molecule 3: DNA (5'-D(*GP*AP*A*CP*TP*CP*AP*AP*CP*AP*GP*AP*CP*CP*GP*TP*GP*CP*CP*GP*CP*CP*TP*GP*CP*CP*T)-3')

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	149.81Å 149.81Å 70.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.25 – 3.99 49.04 – 3.99	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.25-3.99) 99.4 (49.04-3.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.15 (at 4.00Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.217 , 0.282 0.217 , 0.281	Depositor DCC
R_{free} test set	1153 reflections (7.71%)	wwPDB-VP
Wilson B-factor (Å^2)	175.4	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 203.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.054 for -h,-k,l 0.418 for h,-h-k,-l 0.056 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6885	wwPDB-VP
Average B, all atoms (Å^2)	197.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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	0.23	0/1505	0.40	0/2043
1	B	0.23	0/1545	0.41	0/2104
1	C	0.23	0/1559	0.40	0/2111
1	D	0.23	0/1519	0.40	0/2056
2	E	0.52	0/465	0.98	0/718
3	F	0.57	0/493	0.90	1/757 (0.1%)
All	All	0.29	0/7086	0.52	1/9789 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	F	4	DC	O4'-C1'-N1	5.67	111.97	108.00

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	1480	0	1469	20	0
1	B	1518	0	1464	19	0
1	C	1535	0	1542	31	0
1	D	1497	0	1509	33	0
2	E	414	0	226	21	0
3	F	441	0	247	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6885	0	6457	106	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 8.

All (106) 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:43:LYS:HE2	2:E:20:DT:C7	2.02	0.89
1:D:43:LYS:HE2	2:E:20:DT:O4	1.81	0.78
1:D:43:LYS:HE2	2:E:20:DT:H71	1.66	0.75
2:E:11:DC:N4	3:F:17:DG:O6	2.19	0.73
1:C:7:ALA:HB1	1:C:49:HIS:HE1	1.56	0.70
1:D:177:GLU:H	1:D:178:PRO:HD2	1.60	0.67
1:B:19:ALA:HB2	1:B:60:ILE:HD11	1.77	0.66
1:D:159:CYS:O	1:D:161:VAL:N	2.28	0.65
2:E:12:DA:H2"	2:E:13:DC:H5"	1.77	0.65
1:D:136:LEU:HD11	1:D:150:VAL:HG21	1.79	0.65
1:B:172:GLU:OE1	1:B:180:ARG:NH2	2.30	0.65
1:C:28:GLU:HB3	1:D:28:GLU:HB2	1.81	0.61
1:C:6:ARG:N	3:F:24:DG:OP1	2.30	0.61
1:B:72:ALA:HA	1:B:89:LEU:HD21	1.82	0.61
1:A:189:ILE:HG22	1:A:198:ARG:HG3	1.82	0.60
1:D:90:THR:HG21	1:D:185:TRP:HE1	1.66	0.59
1:D:43:LYS:CE	2:E:20:DT:C7	2.78	0.59
1:D:84:GLU:OE2	1:D:88:ARG:NH1	2.36	0.59
1:D:43:LYS:HE2	2:E:20:DT:C4	2.37	0.59
1:D:137:ASP:OD1	1:D:140:ARG:NH1	2.36	0.58
2:E:19:DG:H2"	2:E:20:DT:H5"	1.85	0.58
1:C:159:CYS:SG	1:D:160:SER:N	2.76	0.58
1:A:159:CYS:HA	1:A:162:VAL:HB	1.86	0.57
1:D:43:LYS:CE	2:E:20:DT:O4	2.53	0.57
1:C:182:ALA:HB2	1:C:209:GLU:HG3	1.86	0.56
1:C:26:GLY:O	1:C:30:THR:OG1	2.23	0.56
1:C:7:ALA:HB1	1:C:49:HIS:CE1	2.41	0.55
1:D:69:ARG:NH2	1:D:126:GLU:OE2	2.40	0.55
1:D:14:ILE:HG21	1:D:46:LEU:HG	1.88	0.55
1:A:33:SER:OG	3:F:17:DG:OP2	2.23	0.55
1:B:163:GLY:O	1:B:165:ARG:N	2.41	0.54
1:B:103:LEU:HD23	1:B:103:LEU:H	1.73	0.54
1:C:164:THR:OG1	1:C:165:ARG:N	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:ARG:HG2	1:C:198:ARG:HD3	1.91	0.53
1:C:44:GLY:HA3	2:E:11:DC:H41	1.73	0.53
1:B:14:ILE:O	1:B:18:ALA:N	2.37	0.53
1:C:139:VAL:HA	1:C:144:VAL:HB	1.91	0.52
1:B:96:LEU:HB3	1:B:103:LEU:HD21	1.91	0.52
1:D:26:GLY:O	1:D:28:GLU:N	2.43	0.52
1:C:96:LEU:HD22	1:C:103:LEU:HD13	1.92	0.52
1:C:149:ASP:OD2	1:D:179:ARG:NH2	2.43	0.52
1:B:11:ARG:HE	1:B:15:ILE:HD11	1.75	0.51
1:A:42:THR:OG1	2:E:6:DG:OP2	2.28	0.51
1:D:43:LYS:CE	2:E:20:DT:H72	2.41	0.50
1:A:151:ASP:N	1:A:152:SER:HB3	2.26	0.50
2:E:18:DT:O2	3:F:11:DG:N2	2.44	0.50
1:D:47:TYR:HE2	2:E:20:DT:H5'	1.77	0.50
1:D:139:VAL:HG22	1:D:144:VAL:HB	1.94	0.50
1:D:66:ARG:HA	1:D:69:ARG:HG2	1.93	0.50
1:C:7:ALA:HB2	2:E:8:DC:H3'	1.94	0.49
1:B:131:ALA:O	1:B:135:LEU:N	2.38	0.49
1:B:75:LEU:HD21	1:B:88:ARG:HG3	1.93	0.49
1:A:99:GLN:HG3	1:A:100:GLY:H	1.77	0.49
1:A:32:LEU:N	3:F:18:DC:OP2	2.32	0.48
1:C:21:LEU:HD13	1:C:25:ARG:HH12	1.79	0.48
1:D:144:VAL:HG13	1:D:192:MET:HG2	1.96	0.48
1:B:80:TYR:OH	1:B:88:ARG:NH2	2.47	0.48
3:F:19:DC:H2'	3:F:20:DG:C8	2.48	0.48
1:D:43:LYS:HE2	2:E:20:DT:C5	2.48	0.47
1:A:162:VAL:O	1:A:164:THR:N	2.47	0.47
1:B:149:ASP:HB2	1:B:153:VAL:HG22	1.97	0.47
1:D:162:VAL:HG12	1:D:166:VAL:HG11	1.97	0.47
1:D:100:GLY:O	1:D:104:ARG:NH2	2.48	0.47
1:C:176:ARG:NH2	1:C:179:ARG:HE	2.13	0.46
1:B:14:ILE:HG22	1:B:39:ALA:HB1	1.97	0.46
1:C:164:THR:HG23	1:C:166:VAL:H	1.80	0.46
1:C:176:ARG:HH22	1:C:179:ARG:HE	1.62	0.46
1:D:145:HIS:ND1	1:D:192:MET:O	2.48	0.46
1:D:32:LEU:H	2:E:19:DG:P	2.38	0.45
1:A:15:ILE:HA	1:A:18:ALA:HB3	1.98	0.45
1:A:91:PHE:HE1	1:A:181:LEU:HB3	1.82	0.45
1:D:23:ASP:OD1	1:D:24:ARG:N	2.45	0.45
1:A:28:GLU:HG3	1:B:28:GLU:HB2	1.99	0.45
1:B:140:ARG:H	1:B:140:ARG:HG2	1.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:ARG:HG2	1:A:204:LEU:HD22	1.99	0.44
1:A:15:ILE:HG13	1:A:15:ILE:H	1.60	0.44
1:C:152:SER:HA	1:D:180:ARG:HH12	1.81	0.44
1:B:176:ARG:HD3	1:B:179:ARG:HB2	2.00	0.44
1:C:44:GLY:CA	2:E:11:DC:H41	2.31	0.43
1:A:153:VAL:HB	1:A:154:ALA:H	1.66	0.43
2:E:23:DA:H2"	2:E:24:DG:H2'	2.00	0.43
1:A:105:ALA:O	1:A:109:LEU:N	2.46	0.43
1:C:78:ARG:HD3	1:C:80:TYR:HE1	1.83	0.43
1:A:151:ASP:HA	1:A:153:VAL:HG22	2.01	0.43
1:D:183:GLU:HA	1:D:186:TYR:CD2	2.53	0.43
1:A:181:LEU:O	1:A:184:MET:HG2	2.19	0.43
1:B:101:PRO:HA	1:B:104:ARG:NE	2.34	0.42
1:A:139:VAL:HG22	1:A:144:VAL:HB	2.02	0.42
1:C:181:LEU:HA	1:C:184:MET:HE3	2.01	0.42
3:F:9:DC:H2"	3:F:10:DA:O4'	2.18	0.42
1:B:98:VAL:HG22	1:B:174:ALA:HB1	2.01	0.42
1:D:177:GLU:N	1:D:178:PRO:HD2	2.33	0.42
1:C:176:ARG:HH22	1:C:179:ARG:HB3	1.85	0.41
1:C:204:LEU:HD23	1:C:204:LEU:HA	1.92	0.41
1:C:96:LEU:HB2	1:C:103:LEU:HD22	2.03	0.41
1:C:107:LEU:HG	1:C:127:TRP:HH2	1.84	0.41
1:C:209:GLU:HA	1:C:212:THR:HG22	2.02	0.41
1:C:6:ARG:NH2	2:E:9:DG:H5'	2.35	0.41
1:A:6:ARG:H	1:A:6:ARG:HG3	1.74	0.41
1:C:175:GLY:HA3	1:C:176:ARG:HA	1.70	0.41
1:C:176:ARG:HD2	1:C:176:ARG:HA	1.86	0.41
1:D:181:LEU:HA	1:D:184:MET:HB3	2.02	0.41
1:A:83:LEU:HB2	1:A:138:ALA:HB1	2.02	0.41
1:C:90:THR:HG22	1:C:181:LEU:HD21	2.02	0.40
1:B:157:LEU:O	1:B:159:CYS:N	2.55	0.40
2:E:7:DG:H1'	2:E:8:DC:H5"	2.04	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	188/215 (87%)	160 (85%)	22 (12%)	6 (3%)	4	31
1	B	200/215 (93%)	163 (82%)	29 (14%)	8 (4%)	3	26
1	C	194/215 (90%)	157 (81%)	32 (16%)	5 (3%)	5	34
1	D	185/215 (86%)	154 (83%)	22 (12%)	9 (5%)	2	22
All	All	767/860 (89%)	634 (83%)	105 (14%)	28 (4%)	3	28

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	GLY
1	B	76	ASP
1	B	159	CYS
1	B	162	VAL
1	B	164	THR
1	C	76	ASP
1	D	160	SER
1	A	58	HIS
1	A	177	GLU
1	B	39	ALA
1	D	159	CYS
1	B	80	TYR
1	C	122	HIS
1	C	209	GLU
1	D	27	TYR
1	D	86	LEU
1	A	153	VAL
1	C	51	ALA
1	D	196	THR
1	A	191	GLY
1	B	158	VAL
1	D	177	GLU
1	B	92	GLY
1	C	177	GLU
1	D	121	PRO
1	D	189	ILE
1	A	150	VAL
1	D	41	VAL

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	148/171 (86%)	140 (95%)	8 (5%)	22 50
1	B	148/171 (86%)	133 (90%)	15 (10%)	7 29
1	C	153/171 (90%)	139 (91%)	14 (9%)	9 32
1	D	154/171 (90%)	137 (89%)	17 (11%)	6 26
All	All	603/684 (88%)	549 (91%)	54 (9%)	9 34

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	8	GLU
1	A	28	GLU
1	A	41	VAL
1	A	64	GLN
1	A	136	LEU
1	A	153	VAL
1	A	159	CYS
1	B	11	ARG
1	B	21	LEU
1	B	30	THR
1	B	31	THR
1	B	36	VAL
1	B	54	GLU
1	B	58	HIS
1	B	86	LEU
1	B	89	LEU
1	B	103	LEU
1	B	109	LEU
1	B	140	ARG
1	B	153	VAL
1	B	171	LEU
1	B	196	THR
1	C	28	GLU

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Mol	Chain	Res	Type
1	C	31	THR
1	C	32	LEU
1	C	62	GLU
1	C	64	GLN
1	C	70	ARG
1	C	73	LYS
1	C	93	MET
1	C	117	ARG
1	C	157	LEU
1	C	177	GLU
1	C	195	VAL
1	C	208	LEU
1	C	209	GLU
1	D	22	PHE
1	D	32	LEU
1	D	42	THR
1	D	46	LEU
1	D	56	LEU
1	D	75	LEU
1	D	76	ASP
1	D	108	ARG
1	D	114	VAL
1	D	120	LEU
1	D	125	THR
1	D	132	THR
1	D	147	ASP
1	D	158	VAL
1	D	180	ARG
1	D	197	ARG
1	D	211	GLU

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

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, 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	194/215 (90%)	0.42	4 (2%) 63 54	129, 181, 244, 266	0
1	B	204/215 (94%)	0.34	10 (4%) 29 25	151, 202, 263, 329	0
1	C	200/215 (93%)	0.33	10 (5%) 28 25	147, 201, 264, 281	0
1	D	193/215 (89%)	0.50	17 (8%) 10 9	141, 190, 235, 269	0
2	E	20/27 (74%)	-0.21	0 100 100	161, 186, 281, 299	0
3	F	22/27 (81%)	0.07	1 (4%) 33 27	161, 188, 286, 323	1 (4%)
All	All	833/914 (91%)	0.37	42 (5%) 28 25	129, 193, 258, 329	1 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	164	THR	5.8
3	F	25	DC	5.6
1	B	115	PRO	5.0
1	C	107	LEU	4.9
1	B	186	TYR	4.8
1	D	147	ASP	4.0
1	B	16	GLY	3.9
1	D	54	GLU	3.8
1	C	167	VAL	3.8
1	D	14	ILE	3.4
1	C	163	GLY	3.3
1	B	5	LEU	3.1
1	B	185	TRP	3.0
1	B	109	LEU	2.7
1	A	90	THR	2.7
1	B	35	ILE	2.7
1	A	35	ILE	2.7
1	A	53	LYS	2.7
1	D	115	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	46	LEU	2.6
1	D	192	MET	2.5
1	D	166	VAL	2.5
1	A	107	LEU	2.5
1	D	87	MET	2.5
1	D	15	ILE	2.5
1	D	50	PHE	2.5
1	D	53	LYS	2.3
1	C	93	MET	2.3
1	C	177	GLU	2.3
1	C	14	ILE	2.3
1	D	27	TYR	2.3
1	C	158	VAL	2.2
1	D	55	ASP	2.2
1	D	120	LEU	2.2
1	D	195	VAL	2.1
1	C	192	MET	2.1
1	C	186	TYR	2.1
1	B	121	PRO	2.1
1	B	103	LEU	2.1
1	B	182	ALA	2.0
1	D	105	ALA	2.0
1	D	109	LEU	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.