



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:18 pm BST

PDB ID : 3IWZ
Title : The c-di-GMP Responsive Global Regulator CLP Links Cell-Cell Signaling to Virulence Gene Expression in *Xanthomonas campestris*
Authors : Chin, K.H.; Tu, Z.L.; Tseng, Y.H.; Dow, J.M.; Wang, A.H.J.; Chou, S.H.
Deposited on : 2009-09-03
Resolution : 2.30 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

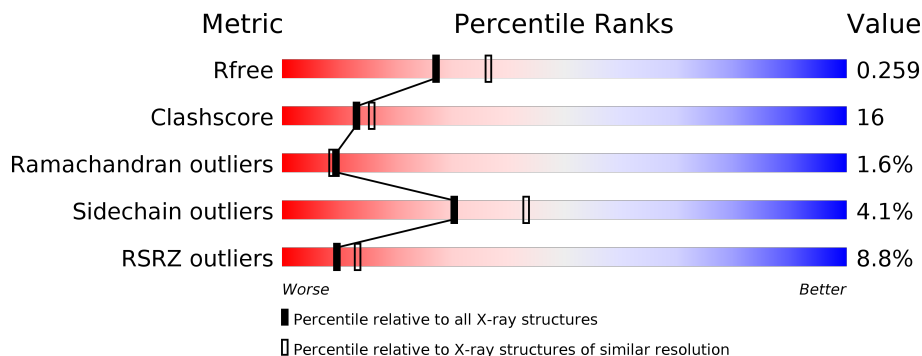
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 $\leq 5\%$. 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	230	 3% 62% 27% 10%
1	B	230	 5% 66% 23% 10%
1	C	230	 15% 54% 30% 5% 10%
1	D	230	 9% 57% 32% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6817 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 Catabolite activation-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	207	Total 1639	C 1025	N 301	O 308	S 5	0	0	0
1	B	207	Total 1639	C 1025	N 301	O 308	S 5	0	0	0
1	C	207	Total 1626	C 1016	N 300	O 305	S 5	0	0	0
1	D	208	Total 1634	C 1022	N 301	O 306	S 5	0	0	0

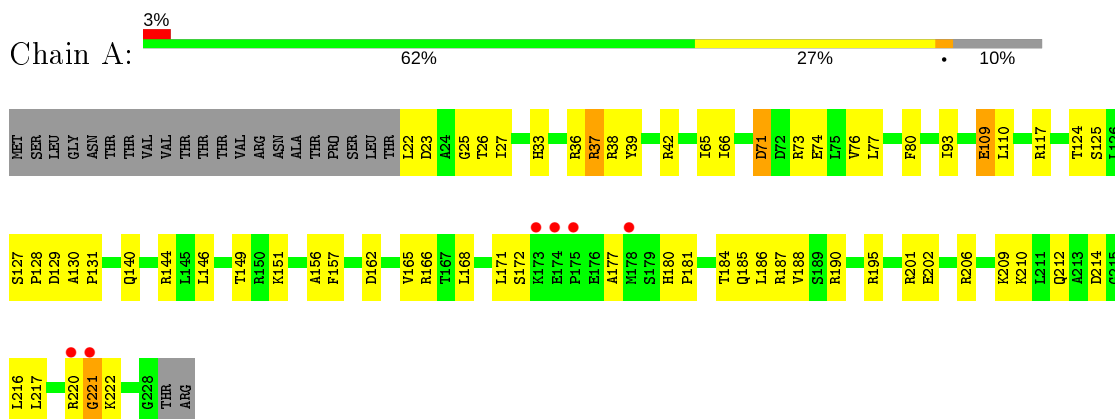
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	96	Total 96	O 96	0	0
2	B	71	Total 71	O 71	0	0
2	C	64	Total 64	O 64	0	0
2	D	48	Total 48	O 48	0	0

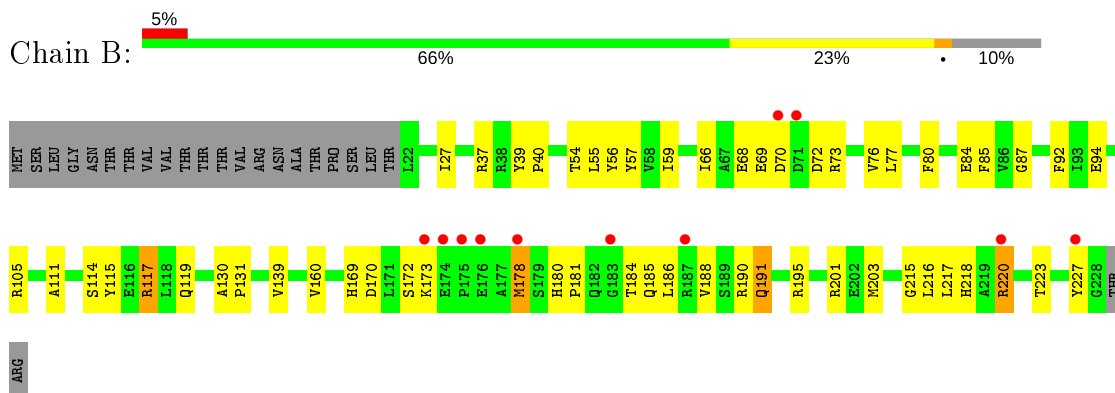
3 Residue-property plots [i](#)

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.

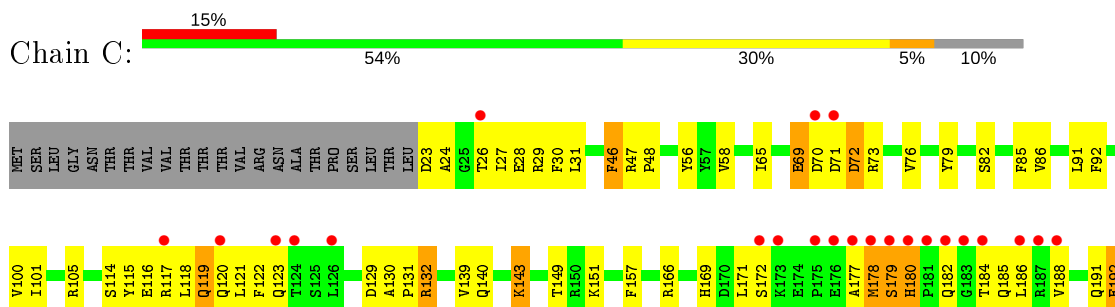
- Molecule 1: Catabolite activation-like protein



- Molecule 1: Catabolite activation-like protein

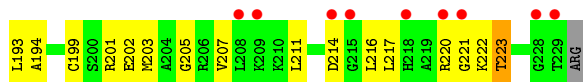
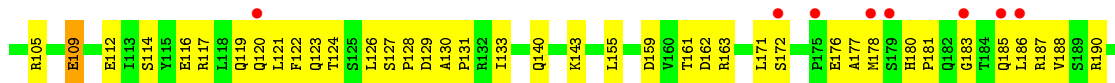


- Molecule 1: Catabolite activation-like protein





• Molecule 1: Catabolite activation-like protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.69Å 67.69Å 110.37Å 90.00° 104.20° 90.00°	Depositor
Resolution (Å)	29.85 – 2.30 29.84 – 2.28	Depositor EDS
% Data completeness (in resolution range)	92.5 (29.85-2.30) 96.9 (29.84-2.28)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.29Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.244 , 0.305 0.257 , 0.259	Depositor DCC
R_{free} test set	2161 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	39.2	Xtrriage
Anisotropy	0.217	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6817	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.35% 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

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	0.38	0/1664	0.62	0/2244
1	B	0.37	0/1664	0.60	0/2244
1	C	0.35	0/1651	0.59	0/2229
1	D	0.33	0/1659	0.59	0/2240
All	All	0.36	0/6638	0.60	0/8957

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	1639	0	1661	51	0
1	B	1639	0	1661	45	0
1	C	1626	0	1638	73	0
1	D	1634	0	1649	58	0
2	A	96	0	0	1	0
2	B	71	0	0	4	0
2	C	64	0	0	2	0
2	D	48	0	0	0	0
All	All	6817	0	6609	214	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 16.

All (214) 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:85:PHE:H	1:D:140:GLN:HE22	1.01	0.99
1:C:85:PHE:H	1:C:140:GLN:HE22	1.05	0.98
1:C:218:HIS:HB3	1:C:225:VAL:HG13	1.49	0.93
1:B:178:MET:HB2	1:B:185:GLN:HB3	1.49	0.93
1:B:185:GLN:HE22	1:B:223:THR:HB	1.34	0.91
1:C:180:HIS:NE2	1:C:182:GLN:HB3	1.92	0.84
1:C:212:GLN:HE22	1:C:218:HIS:HA	1.41	0.84
1:C:222:LYS:HA	1:C:222:LYS:HE2	1.61	0.82
1:D:171:LEU:HD13	1:D:188:VAL:HG21	1.60	0.81
1:C:85:PHE:H	1:C:140:GLN:NE2	1.80	0.79
1:A:76:VAL:HG23	1:A:195:ARG:HB3	1.68	0.75
1:A:37:ARG:NH1	1:A:37:ARG:HB3	2.02	0.74
1:C:130:ALA:HB3	1:C:131:PRO:HD3	1.68	0.74
1:A:130:ALA:HB3	1:A:131:PRO:HD3	1.69	0.73
1:A:117:ARG:HH11	1:A:117:ARG:HG2	1.54	0.72
1:C:116:GLU:O	1:C:120:GLN:HG3	1.90	0.72
1:B:55:LEU:HD23	1:B:56:TYR:N	2.06	0.71
1:D:185:GLN:NE2	1:D:223:THR:HB	2.05	0.71
1:D:121:LEU:HB3	1:D:126:LEU:HD22	1.72	0.70
1:D:129:ASP:O	1:D:133:ILE:HG12	1.90	0.70
1:B:185:GLN:NE2	1:B:223:THR:HB	2.06	0.69
1:D:172:SER:HA	1:D:186:LEU:HD21	1.72	0.69
1:B:130:ALA:HB3	1:B:131:PRO:HD3	1.74	0.69
1:A:71:ASP:OD1	1:A:73:ARG:HB2	1.93	0.68
1:C:30:PHE:HD2	1:C:31:LEU:HD22	1.56	0.68
1:C:85:PHE:N	1:C:140:GLN:HE22	1.87	0.67
1:B:190:ARG:HD3	1:B:201:ARG:HH21	1.59	0.66
1:C:117:ARG:HH22	1:C:121:LEU:HD21	1.61	0.66
1:C:180:HIS:CD2	1:C:182:GLN:HB3	2.30	0.66
1:D:185:GLN:HE22	1:D:223:THR:HB	1.61	0.66
1:C:171:LEU:HD13	1:C:188:VAL:HG11	1.78	0.65
1:B:55:LEU:HD21	1:B:85:PHE:HD1	1.61	0.65
1:D:180:HIS:ND1	1:D:181:PRO:HD2	2.12	0.65
1:D:24:ALA:O	1:D:28:GLU:HG2	1.98	0.64
1:D:70:ASP:C	1:D:72:ASP:H	1.99	0.63
1:A:217:LEU:HD12	1:A:217:LEU:C	2.18	0.63
1:C:218:HIS:HB3	1:C:225:VAL:CG1	2.26	0.63
1:D:85:PHE:H	1:D:140:GLN:NE2	1.84	0.62
1:D:130:ALA:HB3	1:D:131:PRO:HD3	1.80	0.62
1:B:160:VAL:HG21	1:B:203:MET:SD	2.39	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:SER:HB3	1:C:117:ARG:HB2	1.81	0.62
1:A:190:ARG:HD2	1:A:201:ARG:NH2	2.15	0.61
1:B:172:SER:OG	1:B:186:LEU:HD21	2.00	0.61
1:A:80:PHE:CE1	1:A:144:ARG:HD2	2.35	0.61
1:A:131:PRO:HG3	1:D:131:PRO:HG3	1.82	0.61
1:A:146:LEU:HD11	1:D:99:GLU:HG2	1.81	0.61
1:A:149:THR:HG23	1:D:77:LEU:HD11	1.83	0.61
1:D:116:GLU:O	1:D:120:GLN:HG3	2.01	0.61
1:D:194:ALA:HB1	1:D:199:CYS:O	2.01	0.61
1:B:80:PHE:HB3	1:B:84:GLU:HG3	1.83	0.60
1:C:191:GLN:HE22	1:C:201:ARG:HD2	1.66	0.60
1:A:171:LEU:HD13	1:A:188:VAL:HG21	1.84	0.59
1:D:119:GLN:HA	1:D:122:PHE:HB2	1.82	0.59
1:D:80:PHE:HB3	1:D:84:GLU:HG3	1.84	0.59
1:A:127:SER:HB3	1:A:128:PRO:HD3	1.84	0.59
1:B:115:TYR:O	1:B:119:GLN:HG2	2.03	0.59
1:D:187:ARG:HA	1:D:223:THR:HG22	1.85	0.58
1:C:220:ARG:HG2	1:C:220:ARG:HH11	1.69	0.58
1:A:117:ARG:HG2	1:A:117:ARG:NH1	2.19	0.58
1:B:69:GLU:OE1	1:B:69:GLU:HA	2.03	0.57
1:C:191:GLN:NE2	1:C:201:ARG:HD2	2.19	0.57
1:B:114:SER:OG	1:B:117:ARG:HB2	2.04	0.57
1:A:212:GLN:HA	1:A:217:LEU:O	2.04	0.57
1:A:80:PHE:HE1	1:A:144:ARG:HD2	1.70	0.57
1:C:218:HIS:HD2	1:C:225:VAL:HG11	1.69	0.57
1:D:190:ARG:HH11	1:D:205:GLY:HA3	1.69	0.56
1:B:178:MET:CB	1:B:185:GLN:HB3	2.29	0.56
1:C:91:LEU:HD11	1:C:118:LEU:HD23	1.88	0.56
1:D:172:SER:OG	1:D:186:LEU:HD11	2.05	0.56
1:B:105:ARG:HD3	2:B:258:HOH:O	2.05	0.56
1:D:207:VAL:O	1:D:211:LEU:HG	2.05	0.56
1:C:24:ALA:O	1:C:28:GLU:HG3	2.06	0.55
1:C:47:ARG:HB3	1:C:47:ARG:NH1	2.21	0.55
1:C:79:TYR:CD1	1:C:105:ARG:HD2	2.41	0.55
1:D:217:LEU:C	1:D:217:LEU:HD12	2.26	0.55
1:A:172:SER:HB3	1:A:186:LEU:HD11	1.88	0.54
1:D:85:PHE:N	1:D:140:GLN:HE22	1.86	0.54
1:D:161:THR:HG22	1:D:207:VAL:HG22	1.88	0.54
1:C:132:ARG:HG3	2:C:241:HOH:O	2.05	0.54
1:C:23:ASP:O	1:C:27:ILE:HG12	2.08	0.54
1:D:79:TYR:CD1	1:D:105:ARG:HD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:221:GLY:C	1:C:223:THR:H	2.12	0.53
1:A:220:ARG:HH22	1:C:166:ARG:HH22	1.54	0.53
1:B:218:HIS:CD2	1:B:220:ARG:HB2	2.43	0.53
1:A:26:THR:HG23	1:A:129:ASP:HB3	1.91	0.53
1:B:92:PHE:HB3	1:C:139:VAL:HA	1.91	0.53
1:C:30:PHE:CD2	1:C:31:LEU:HD22	2.41	0.53
1:C:169:HIS:ND1	1:C:226:LEU:HD21	2.24	0.53
1:A:210:LYS:HE3	1:A:214:ASP:OD1	2.10	0.52
1:A:93:ILE:HD11	1:D:143:LYS:HG3	1.91	0.52
1:B:59:ILE:HD11	1:B:111:ALA:HB2	1.92	0.52
1:A:37:ARG:CZ	1:A:37:ARG:HB3	2.39	0.52
1:A:220:ARG:NH2	1:C:166:ARG:HH22	2.08	0.52
1:C:194:ALA:HB1	1:C:199:CYS:O	2.09	0.52
1:C:69:GLU:CD	1:C:70:ASP:H	2.14	0.52
1:C:172:SER:OG	1:C:186:LEU:HD21	2.10	0.51
1:B:172:SER:HA	1:B:186:LEU:HD21	1.92	0.51
1:D:70:ASP:C	1:D:72:ASP:N	2.64	0.51
1:C:201:ARG:HH22	1:C:222:LYS:HZ1	1.57	0.51
1:B:188:VAL:HG12	2:B:275:HOH:O	2.09	0.51
1:C:132:ARG:HE	1:C:132:ARG:HA	1.76	0.51
1:B:172:SER:OG	1:B:186:LEU:HD11	2.11	0.51
1:C:65:ILE:HG23	1:C:100:VAL:HG21	1.93	0.51
1:A:157:PHE:HB2	1:D:69:GLU:HG3	1.93	0.51
1:A:39:TYR:CE2	1:A:110:LEU:HD12	2.46	0.50
1:D:56:TYR:HB2	1:D:86:VAL:HG23	1.93	0.50
1:D:22:LEU:O	1:D:22:LEU:HD23	2.12	0.50
1:C:186:LEU:HB2	1:C:224:VAL:HB	1.93	0.50
1:C:218:HIS:CD2	1:C:225:VAL:HG11	2.47	0.50
1:C:212:GLN:HB2	1:C:217:LEU:HD11	1.93	0.50
1:D:127:SER:HB3	1:D:128:PRO:HD3	1.93	0.50
1:A:140:GLN:O	1:A:144:ARG:HD3	2.12	0.49
1:B:68:GLU:HG2	1:B:69:GLU:N	2.27	0.49
1:B:191:GLN:NE2	1:B:201:ARG:HD2	2.27	0.49
1:A:156:ALA:HB2	1:D:155:LEU:HD13	1.94	0.49
1:D:171:LEU:CD1	1:D:188:VAL:HG21	2.38	0.49
1:B:37:ARG:HH11	1:B:37:ARG:HG2	1.77	0.49
1:D:201:ARG:HG3	1:D:201:ARG:HH11	1.77	0.49
1:D:70:ASP:O	1:D:72:ASP:N	2.45	0.49
1:C:178:MET:O	1:C:179:SER:HB2	2.13	0.49
1:A:221:GLY:HA3	1:B:70:ASP:O	2.13	0.48
1:A:214:ASP:HB3	1:A:216:LEU:HD13	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:LEU:O	1:D:217:LEU:HD12	2.14	0.48
1:B:217:LEU:C	1:B:217:LEU:HD12	2.35	0.48
1:C:192:GLU:OE1	1:C:192:GLU:HA	2.14	0.48
1:C:27:ILE:O	1:C:31:LEU:HD23	2.14	0.48
1:C:29:ARG:CZ	1:C:29:ARG:HB2	2.44	0.48
1:A:180:HIS:HB2	1:A:185:GLN:HB2	1.96	0.47
1:C:71:ASP:O	1:C:72:ASP:CB	2.61	0.47
1:A:165:VAL:HG23	1:A:166:ARG:N	2.30	0.47
1:D:123:GLN:HA	1:D:123:GLN:OE1	2.14	0.47
1:D:220:ARG:O	1:D:222:LYS:N	2.47	0.47
1:C:91:LEU:HG	1:C:115:TYR:CD1	2.50	0.47
1:C:119:GLN:O	1:C:123:GLN:HG3	2.14	0.47
1:A:37:ARG:HB3	1:A:37:ARG:HH11	1.77	0.47
1:A:38:ARG:NH1	2:A:296:HOH:O	2.47	0.47
1:A:177:ALA:HB1	1:A:184:THR:CG2	2.45	0.47
1:A:168:LEU:O	1:A:186:LEU:HD21	2.14	0.47
1:A:210:LYS:O	1:A:210:LYS:HD3	2.15	0.46
1:C:71:ASP:HB3	1:C:73:ARG:HG2	1.97	0.46
1:A:187:ARG:HD3	1:A:222:LYS:O	2.16	0.46
1:A:65:ILE:HG22	1:A:77:LEU:HD12	1.98	0.46
1:A:124:THR:OG1	1:A:125:SER:N	2.47	0.46
1:D:71:ASP:N	1:D:71:ASP:OD2	2.49	0.46
1:C:58:VAL:HG12	1:C:82:SER:HA	1.98	0.45
1:D:114:SER:OG	1:D:117:ARG:HB2	2.16	0.45
1:A:76:VAL:O	1:A:151:LYS:HE3	2.16	0.45
1:C:212:GLN:NE2	1:C:219:ALA:H	2.14	0.45
1:A:190:ARG:HD2	1:A:201:ARG:HH21	1.81	0.45
1:B:139:VAL:HA	1:C:92:PHE:HB3	1.99	0.45
1:C:184:THR:OG1	1:C:226:LEU:HB3	2.16	0.45
1:B:55:LEU:C	1:B:55:LEU:HD23	2.36	0.45
1:C:220:ARG:HH11	1:C:220:ARG:CG	2.28	0.45
1:C:47:ARG:CB	1:C:47:ARG:HH11	2.29	0.45
1:D:187:ARG:CA	1:D:223:THR:HG22	2.46	0.45
1:D:121:LEU:O	1:D:126:LEU:HB2	2.17	0.45
1:B:227:TYR:CD2	1:B:227:TYR:N	2.84	0.44
1:C:48:PRO:HD3	1:C:101:ILE:HG12	1.98	0.44
1:B:169:HIS:O	1:B:173:LYS:HG2	2.17	0.44
1:B:77:LEU:HD11	1:C:149:THR:HG23	2.00	0.44
1:B:87:GLY:HA2	2:B:240:HOH:O	2.17	0.44
1:B:172:SER:CA	1:B:186:LEU:HD21	2.48	0.44
1:D:159:ASP:O	1:D:162:ASP:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:VAL:HG23	1:B:195:ARG:HB3	1.99	0.44
1:D:117:ARG:O	1:D:121:LEU:HG	2.18	0.44
1:B:66:ILE:HG12	1:B:76:VAL:HG22	1.99	0.44
1:B:180:HIS:ND1	1:B:181:PRO:HD2	2.33	0.43
1:B:215:GLY:O	1:B:216:LEU:HD23	2.18	0.43
1:C:211:LEU:O	1:C:214:ASP:HB2	2.18	0.43
1:C:71:ASP:O	1:C:72:ASP:HB3	2.18	0.43
1:B:220:ARG:HH11	1:B:220:ARG:HG2	1.83	0.43
1:D:121:LEU:HD13	1:D:126:LEU:HD22	2.00	0.43
1:A:210:LYS:HD3	1:A:210:LYS:C	2.38	0.43
1:A:33:HIS:HD2	1:A:117:ARG:NH2	2.17	0.43
1:A:171:LEU:HD13	1:A:188:VAL:CG2	2.48	0.43
1:B:94:GLU:HA	2:B:267:HOH:O	2.18	0.43
1:A:162:ASP:HA	1:A:165:VAL:HG22	2.00	0.43
1:D:159:ASP:O	1:D:163:ARG:HG3	2.19	0.43
1:B:27:ILE:HD13	1:B:57:TYR:OH	2.19	0.43
1:D:117:ARG:NH2	1:D:121:LEU:HD21	2.34	0.43
1:D:190:ARG:HD2	1:D:205:GLY:CA	2.49	0.43
1:C:26:THR:HG23	1:C:129:ASP:HB3	1.99	0.43
1:C:46:PHE:C	1:C:46:PHE:CD1	2.92	0.43
1:C:76:VAL:HG23	1:C:195:ARG:HB3	2.00	0.42
1:D:55:LEU:HD22	1:D:56:TYR:N	2.34	0.42
1:A:22:LEU:HG	1:A:27:ILE:HG13	2.01	0.42
1:A:190:ARG:HE	1:A:201:ARG:HE	1.67	0.42
1:A:66:ILE:CG2	1:A:74:GLU:HB3	2.49	0.42
1:C:201:ARG:NH2	1:C:222:LYS:NZ	2.67	0.42
1:C:114:SER:HB3	1:C:117:ARG:CB	2.48	0.42
1:A:36:ARG:HB3	1:A:109:GLU:OE2	2.20	0.42
1:B:54:THR:HA	1:B:114:SER:HA	2.02	0.41
1:C:119:GLN:HA	1:C:122:PHE:HB2	2.02	0.41
1:D:117:ARG:HA	1:D:117:ARG:HD2	1.84	0.41
1:C:132:ARG:CA	1:C:132:ARG:HE	2.33	0.41
1:C:76:VAL:O	1:C:151:LYS:HE2	2.20	0.41
1:C:56:TYR:HB2	1:C:86:VAL:HG23	2.03	0.41
1:D:66:ILE:HB	1:D:101:ILE:HB	2.03	0.41
1:C:208:LEU:O	1:C:212:GLN:HB2	2.20	0.41
1:C:178:MET:HB2	1:C:185:GLN:H	1.86	0.41
1:B:39:TYR:HA	1:B:40:PRO:HD3	1.92	0.41
1:B:57:TYR:HB3	1:B:111:ALA:HB3	2.02	0.41
1:A:23:ASP:OD1	1:A:25:GLY:N	2.53	0.41
1:D:193:LEU:HD12	1:D:193:LEU:HA	1.95	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:54:THR:HG21	1:D:112:GLU:OE2	2.21	0.41
1:C:222:LYS:CE	1:C:222:LYS:HA	2.42	0.41
1:A:180:HIS:ND1	1:A:181:PRO:N	2.69	0.40
1:C:201:ARG:HH22	1:C:222:LYS:NZ	2.19	0.40
1:D:93:ILE:O	1:D:94:GLU:C	2.59	0.40
1:C:143:LYS:HD3	2:C:284:HOH:O	2.20	0.40
1:D:38:ARG:CD	1:D:109:GLU:HG3	2.51	0.40
1:B:73:ARG:CZ	1:C:157:PHE:HB3	2.51	0.40
1:B:191:GLN:HE21	1:B:191:GLN:HB2	1.57	0.40
1:D:176:GLU:O	1:D:177:ALA:C	2.60	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	205/230 (89%)	198 (97%)	6 (3%)	1 (0%)	29	35
1	B	205/230 (89%)	198 (97%)	6 (3%)	1 (0%)	29	35
1	C	205/230 (89%)	185 (90%)	14 (7%)	6 (3%)	4	3
1	D	206/230 (90%)	191 (93%)	10 (5%)	5 (2%)	6	4
All	All	821/920 (89%)	772 (94%)	36 (4%)	13 (2%)	9	9

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	179	SER
1	D	221	GLY
1	D	71	ASP
1	A	221	GLY
1	C	178	MET

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Mol	Chain	Res	Type
1	C	72	ASP
1	C	180	HIS
1	C	222	LYS
1	D	216	LEU
1	D	214	ASP
1	B	178	MET
1	C	177	ALA
1	D	183	GLY

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	177/198 (89%)	170 (96%)	7 (4%)	31	44
1	B	177/198 (89%)	171 (97%)	6 (3%)	37	51
1	C	174/198 (88%)	167 (96%)	7 (4%)	31	44
1	D	175/198 (88%)	166 (95%)	9 (5%)	24	33
All	All	703/792 (89%)	674 (96%)	29 (4%)	30	43

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	42	ARG
1	A	71	ASP
1	A	109	GLU
1	A	202	GLU
1	A	206	ARG
1	A	209	LYS
1	B	72	ASP
1	B	117	ARG
1	B	170	ASP
1	B	184	THR
1	B	191	GLN
1	B	220	ARG

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Mol	Chain	Res	Type
1	C	46	PHE
1	C	69	GLU
1	C	119	GLN
1	C	132	ARG
1	C	143	LYS
1	C	192	GLU
1	C	220	ARG
1	D	22	LEU
1	D	36	ARG
1	D	55	LEU
1	D	109	GLU
1	D	124	THR
1	D	178	MET
1	D	202	GLU
1	D	203	MET
1	D	223	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	107	GLN
1	A	169	HIS
1	B	120	GLN
1	B	123	GLN
1	B	185	GLN
1	B	191	GLN
1	C	33	HIS
1	C	120	GLN
1	C	140	GLN
1	C	191	GLN
1	C	212	GLN
1	C	218	HIS
1	D	120	GLN
1	D	140	GLN
1	D	185	GLN
1	D	212	GLN

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](#)

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	207/230 (90%)	0.20	6 (2%) 51 58	21, 39, 59, 63	0
1	B	207/230 (90%)	0.47	11 (5%) 26 33	22, 40, 63, 68	0
1	C	207/230 (90%)	1.02	35 (16%) 1 2	22, 49, 68, 72	0
1	D	208/230 (90%)	0.69	21 (10%) 7 9	22, 49, 65, 69	0
All	All	829/920 (90%)	0.59	73 (8%) 10 13	21, 43, 65, 72	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	229	THR	9.5
1	D	229	THR	8.3
1	D	183	GLY	7.9
1	C	183	GLY	7.5
1	C	226	LEU	6.8
1	C	219	ALA	6.2
1	C	228	GLY	5.9
1	C	186	LEU	5.4
1	C	181	PRO	5.3
1	C	217	LEU	5.2
1	B	178	MET	5.1
1	B	220	ARG	5.1
1	C	178	MET	5.0
1	C	180	HIS	4.6
1	D	178	MET	4.5
1	C	175	PRO	4.5
1	C	177	ALA	4.4
1	B	174	GLU	4.4
1	B	175	PRO	4.3
1	D	228	GLY	4.3
1	C	179	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	184	THR	3.8
1	C	220	ARG	3.8
1	B	183	GLY	3.7
1	D	186	LEU	3.7
1	D	70	ASP	3.7
1	C	218	HIS	3.6
1	C	187	ARG	3.5
1	B	187	ARG	3.5
1	D	179	SER	3.4
1	C	227	TYR	3.3
1	D	220	ARG	3.1
1	D	120	GLN	3.1
1	D	208	LEU	3.0
1	C	188	VAL	3.0
1	B	227	TYR	2.9
1	D	218	HIS	2.9
1	C	225	VAL	2.9
1	D	221	GLY	2.8
1	C	123	GLN	2.8
1	A	173	LYS	2.8
1	D	35	HIS	2.8
1	D	185	GLN	2.8
1	C	124	THR	2.7
1	C	182	GLN	2.7
1	D	209	LYS	2.7
1	A	221	GLY	2.6
1	A	174	GLU	2.6
1	B	70	ASP	2.6
1	D	32	ALA	2.6
1	C	173	LYS	2.6
1	D	172	SER	2.5
1	C	126	LEU	2.5
1	C	70	ASP	2.5
1	C	213	ALA	2.5
1	D	71	ASP	2.5
1	C	71	ASP	2.5
1	C	117	ARG	2.4
1	C	176	GLU	2.3
1	D	214	ASP	2.3
1	B	173	LYS	2.3
1	C	209	LYS	2.3
1	D	175	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	215	GLY	2.2
1	C	120	GLN	2.2
1	B	71	ASP	2.2
1	A	178	MET	2.1
1	C	172	SER	2.1
1	B	176	GLU	2.1
1	C	224	VAL	2.1
1	A	175	PRO	2.0
1	C	26	THR	2.0
1	A	220	ARG	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.