



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:38 pm BST

PDB ID : 3Q4J
Title : Structure of a small peptide ligand bound to E.coli DNA sliding clamp
Authors : Wolff, P.; Olieric, V.; Briand, J.P.; Chaloin, O.; Dejaegere, A.; Dumas, P.;
Ennifar, E.; Guichard, G.; Wagner, J.; Burnouf, D.
Deposited on : 2010-12-23
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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

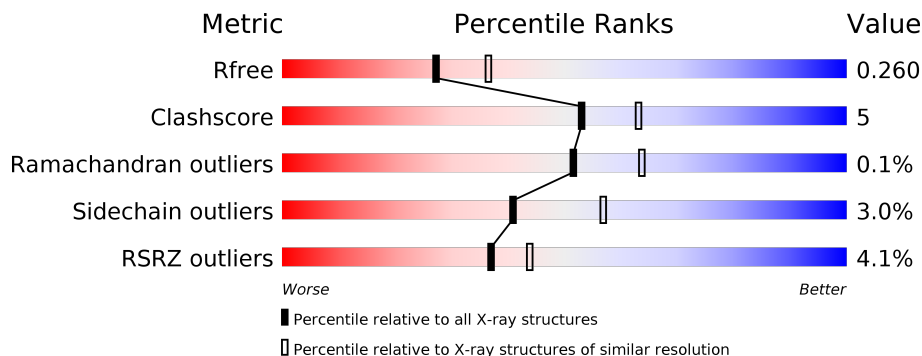
1 Overall quality at a glance

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	366	
1	B	366	
1	C	366	
1	D	366	
1	E	366	
1	F	366	

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Mol	Chain	Length	Quality of chain
2	H	6	 67% 33%
2	I	6	 50% 83% 17%
2	J	6	 50% 67% 33%
2	K	6	 67% 33%
2	L	6	 50% 50%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 17322 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 DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	366	Total 2774	C 1747	N 479	O 530	S 18	0	1	0
1	B	364	Total 2792	C 1756	N 485	O 532	S 19	0	0	0
1	C	365	Total 2772	C 1748	N 472	O 533	S 19	0	0	0
1	D	365	Total 2779	C 1748	N 482	O 530	S 19	0	0	0
1	E	364	Total 2777	C 1744	N 482	O 532	S 19	0	0	0
1	F	365	Total 2800	C 1762	N 486	O 533	S 19	0	0	0

- Molecule 2 is a protein called peptide ligand.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	H	6	Total 48	C 32	N 6	O 10	0	0	0
2	I	6	Total 48	C 32	N 6	O 10	0	0	0
2	J	6	Total 48	C 32	N 6	O 10	0	0	0
2	K	6	Total 48	C 32	N 6	O 10	0	0	0
2	L	6	Total 48	C 32	N 6	O 10	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total 48	O 48	0	0

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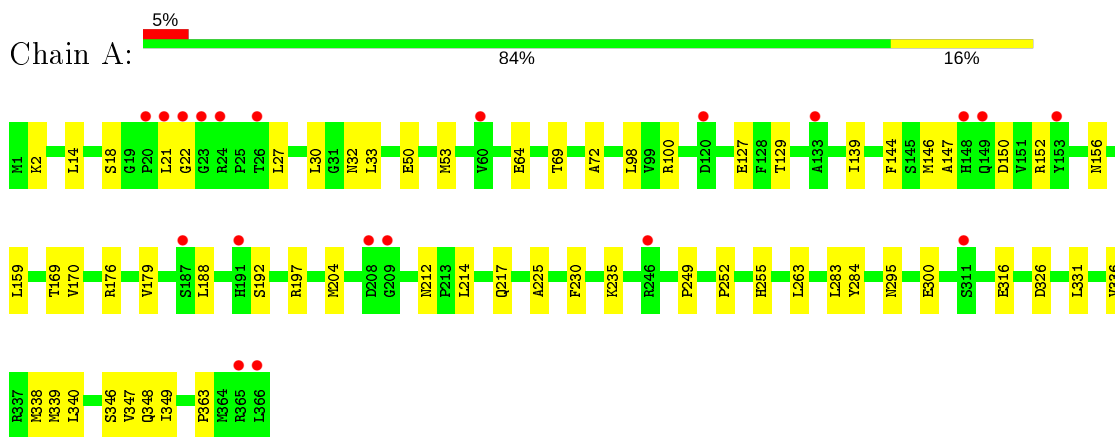
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	98	Total O 98 98	0	0
3	C	44	Total O 44 44	0	0
3	D	46	Total O 46 46	0	0
3	E	75	Total O 75 75	0	0
3	F	72	Total O 72 72	0	0
3	H	3	Total O 3 3	0	0
3	K	2	Total O 2 2	0	0

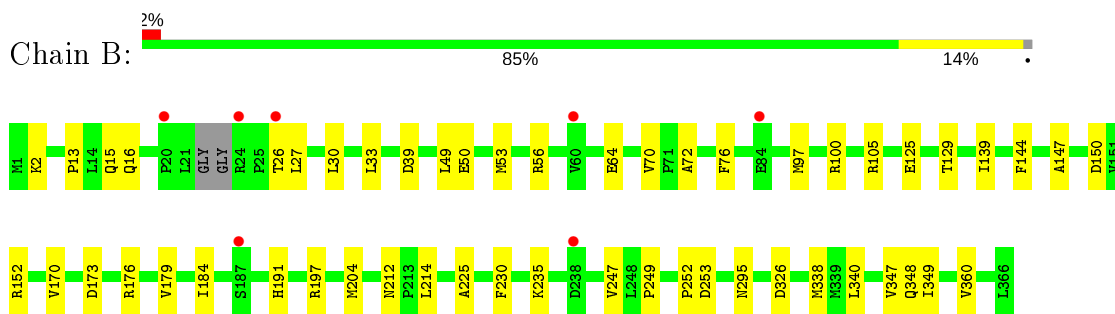
3 Residue-property plots [i](#)

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.

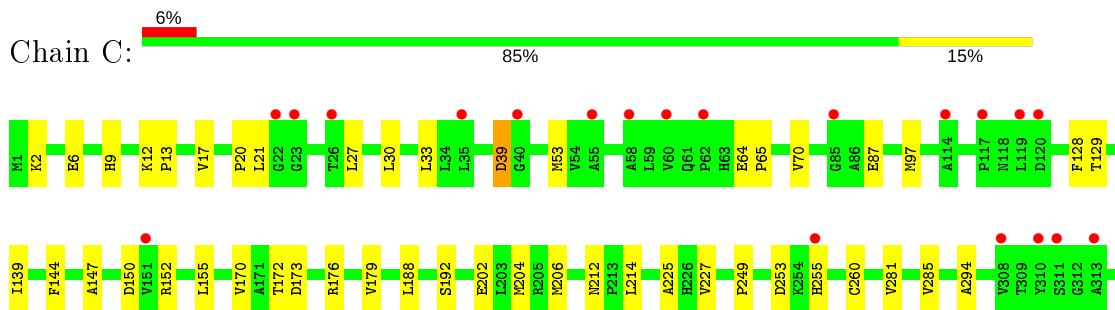
- Molecule 1: DNA polymerase III subunit beta

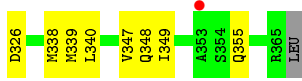


- Molecule 1: DNA polymerase III subunit beta

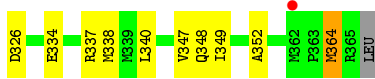
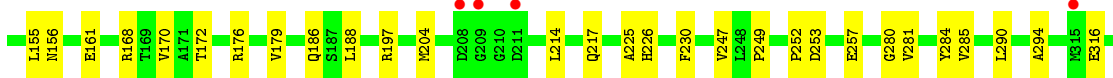
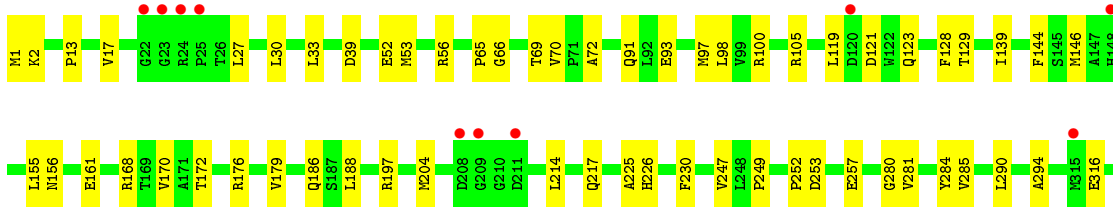
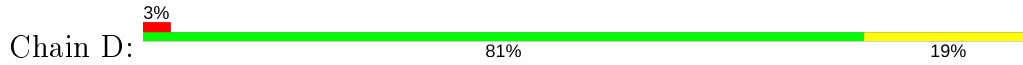


- Molecule 1: DNA polymerase III subunit beta

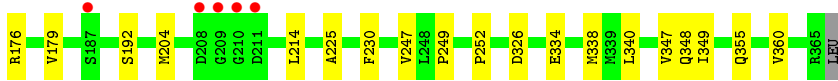
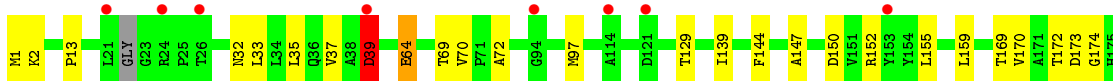
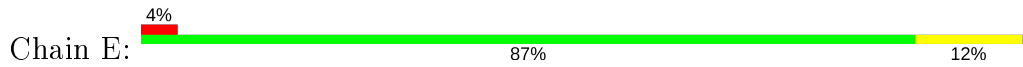




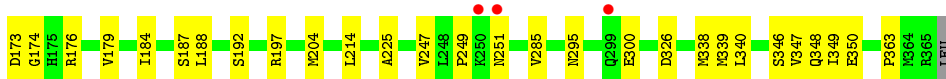
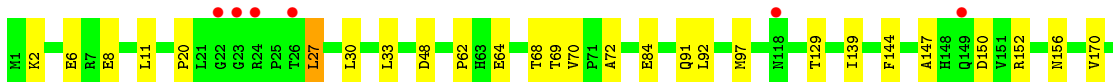
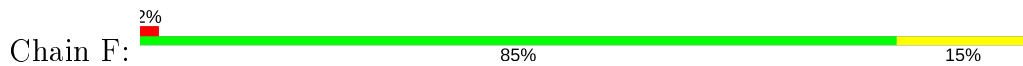
- Molecule 1: DNA polymerase III subunit beta



- Molecule 1: DNA polymerase III subunit beta



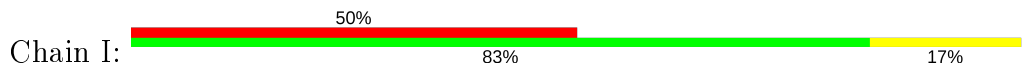
- Molecule 1: DNA polymerase III subunit beta



- Molecule 2: peptide ligand



- Molecule 2: peptide ligand





- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



- Molecule 2: peptide ligand



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	35.09Å 132.87Å 137.27Å 62.73° 88.51° 89.77°	Depositor
Resolution (Å)	29.52 – 2.30 29.52 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.52-2.30) 98.6 (29.52-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.31Å)	Xtrriage
Refinement program	BUSTER-TNT BUSTER 2.8.0, BUSTER 2.8.0	Depositor
R, R_{free}	0.214 , 0.254 0.220 , 0.260	Depositor DCC
R_{free} test set	1959 reflections (2.03%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtrriage
Anisotropy	0.487	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-l 0.000 for -h,k,k-l 0.000 for -h,-k,-k+l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17322	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.06% 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: ACE

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	0.46	0/2827	0.69	0/3841
1	B	0.53	0/2840	0.70	0/3851
1	C	0.49	0/2821	0.73	0/3829
1	D	0.48	0/2826	0.70	0/3837
1	E	0.51	0/2825	0.70	1/3831 (0.0%)
1	F	0.53	0/2847	0.72	0/3860
2	H	0.69	0/46	0.70	0/60
2	I	0.58	0/46	0.67	0/60
2	J	0.55	0/46	0.67	0/60
2	K	0.57	0/46	0.65	0/60
2	L	0.56	0/46	0.61	0/60
All	All	0.50	0/17216	0.70	1/23349 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	E	39	ASP	N-CA-C	5.75	126.51	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2774	0	2722	35	0
1	B	2792	0	2771	27	0
1	C	2772	0	2731	26	0
1	D	2779	0	2738	41	0
1	E	2777	0	2726	27	0
1	F	2800	0	2789	27	0
2	H	48	0	46	2	0
2	I	48	0	46	0	0
2	J	48	0	46	1	0
2	K	48	0	46	3	0
2	L	48	0	46	2	0
3	A	48	0	0	0	0
3	B	98	0	0	0	0
3	C	44	0	0	1	0
3	D	46	0	0	0	0
3	E	75	0	0	0	0
3	F	72	0	0	0	0
3	H	3	0	0	0	0
3	K	2	0	0	0	0
All	All	17322	0	16707	180	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 5.

All (180) 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:A:32:ASN:HB3	1:A:69:THR:HG22	1.28	1.13
1:D:280:GLY:HA2	1:D:364:MET:HE1	1.35	1.09
1:E:32:ASN:HB3	1:E:69:THR:HG22	1.33	1.05
1:D:280:GLY:CA	1:D:364:MET:HE1	1.98	0.91
1:D:1:MET:HB3	1:D:66:GLY:HA3	1.57	0.86
1:C:260:CYS:SG	3:C:415:HOH:O	2.42	0.77
1:A:127:GLU:HB2	1:C:65:PRO:HG3	1.70	0.73
1:A:32:ASN:CB	1:A:69:THR:HG22	2.14	0.73
1:D:280:GLY:HA2	1:D:364:MET:CE	2.15	0.73
1:A:98:LEU:HD23	1:A:100:ARG:HH12	1.50	0.72
1:E:249:PRO:HD2	1:E:348:GLN:HE21	1.54	0.71
1:F:2:LYS:HB2	1:F:64:GLU:HB2	1.73	0.70
1:A:249:PRO:HD2	1:A:348:GLN:HE21	1.56	0.70
1:B:249:PRO:HD2	1:B:348:GLN:HE21	1.57	0.69
1:B:16:GLN:HE21	1:B:53:MET:HE2	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:214:LEU:HD11	1:C:225:ALA:HB1	1.75	0.68
1:C:249:PRO:HD2	1:C:348:GLN:HE21	1.58	0.67
1:E:214:LEU:HD11	1:E:225:ALA:HB1	1.76	0.67
1:E:1:MET:HE2	1:E:37:VAL:HG23	1.77	0.67
1:E:32:ASN:CB	1:E:69:THR:HG22	2.19	0.66
1:D:214:LEU:HD11	1:D:225:ALA:HB1	1.77	0.66
1:D:98:LEU:HD23	1:D:100:ARG:HH12	1.59	0.66
1:A:249:PRO:HB2	1:A:252:PRO:HG3	1.78	0.66
1:F:139:ILE:HG21	1:F:204:MET:HG2	1.76	0.66
1:A:214:LEU:HD11	1:A:225:ALA:HB1	1.78	0.65
1:B:100:ARG:HG2	1:B:105:ARG:HG2	1.78	0.65
1:D:1:MET:HB3	1:D:66:GLY:CA	2.26	0.65
1:C:338:MET:HG2	1:C:349:ILE:HG12	1.79	0.65
1:D:249:PRO:HD2	1:D:348:GLN:HE21	1.62	0.64
1:F:214:LEU:HD11	1:F:225:ALA:HB1	1.80	0.63
1:B:53:MET:HE3	1:B:230:PHE:HB3	1.81	0.62
1:F:249:PRO:HD2	1:F:348:GLN:HE21	1.64	0.62
1:B:53:MET:CE	1:B:230:PHE:HB3	2.30	0.62
1:D:139:ILE:HG21	1:D:204:MET:HG2	1.82	0.61
1:E:139:ILE:HG21	1:E:204:MET:HG2	1.83	0.60
1:B:139:ILE:HG21	1:B:204:MET:HG2	1.84	0.60
1:C:281:VAL:HG12	1:C:294:ALA:HB2	1.84	0.60
1:B:214:LEU:HD11	1:B:225:ALA:HB1	1.82	0.59
1:C:139:ILE:HG21	1:C:204:MET:HG2	1.84	0.59
1:F:20:PRO:O	1:F:48:ASP:HB3	2.03	0.58
1:F:70:VAL:HG11	1:F:97:MET:SD	2.44	0.58
1:E:247:VAL:HG13	2:K:73:LEU:HD22	1.86	0.58
1:D:128:PHE:HA	1:D:186:GLN:HE22	1.71	0.56
1:F:68:THR:HG21	1:F:92:LEU:HD21	1.87	0.56
1:A:139:ILE:HG21	1:A:204:MET:HG2	1.87	0.56
1:F:6:GLU:OE2	1:F:8:GLU:HG2	2.06	0.55
1:F:338:MET:HE3	1:F:347:VAL:HG21	1.89	0.55
1:E:340:LEU:HG	1:E:347:VAL:HG23	1.89	0.55
1:D:65:PRO:HD2	1:F:187:SER:O	2.07	0.55
1:C:20:PRO:HG2	1:C:53:MET:HG2	1.90	0.54
1:E:1:MET:CE	1:E:37:VAL:HG23	2.38	0.54
1:A:284:TYR:CE1	1:A:316:GLU:HG3	2.43	0.54
1:B:170:VAL:HG22	1:B:179:VAL:HG23	1.89	0.54
1:F:247:VAL:HG13	2:L:73:LEU:HD22	1.90	0.54
1:A:32:ASN:HB3	1:A:69:THR:CG2	2.20	0.53
1:C:128:PHE:HB3	1:C:188:LEU:HD21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:338:MET:HG2	1:E:349:ILE:HG12	1.91	0.53
1:E:39:ASP:CG	1:E:39:ASP:O	2.47	0.53
1:E:1:MET:HE1	1:E:35:LEU:HB3	1.91	0.53
1:D:249:PRO:HB2	1:D:252:PRO:HG3	1.91	0.52
1:D:284:TYR:CE1	1:D:316:GLU:HG3	2.44	0.52
1:D:340:LEU:HG	1:D:347:VAL:HG23	1.89	0.52
1:F:340:LEU:HG	1:F:347:VAL:HG23	1.91	0.52
1:A:340:LEU:HG	1:A:347:VAL:HG23	1.91	0.51
1:B:150:ASP:OD2	1:B:152:ARG:HD3	2.11	0.51
1:A:2:LYS:HB3	1:A:64:GLU:HB3	1.93	0.51
1:D:39:ASP:OD2	1:D:39:ASP:CB	2.58	0.51
1:E:249:PRO:HB2	1:E:252:PRO:HG3	1.93	0.51
1:B:249:PRO:HB2	1:B:252:PRO:HG3	1.92	0.51
1:C:340:LEU:HG	1:C:347:VAL:HG23	1.93	0.51
1:E:170:VAL:HG22	1:E:179:VAL:HG23	1.92	0.51
1:B:340:LEU:HG	1:B:347:VAL:HG23	1.92	0.51
1:A:170:VAL:HG22	1:A:179:VAL:HG23	1.92	0.50
1:C:2:LYS:HB3	1:C:64:GLU:HB3	1.92	0.50
1:C:6:GLU:HB3	1:C:9:HIS:HD2	1.76	0.50
1:D:52:GLU:HG2	1:D:119:LEU:HD22	1.93	0.50
1:B:144:PHE:CD2	1:B:326:ASP:HB3	2.47	0.50
1:B:13:PRO:HA	1:B:230:PHE:HE1	1.76	0.50
1:D:13:PRO:HA	1:D:230:PHE:HE1	1.77	0.50
1:D:280:GLY:CA	1:D:364:MET:CE	2.82	0.50
1:D:338:MET:HG2	1:D:349:ILE:HG12	1.93	0.50
1:D:33:LEU:HG	1:D:72:ALA:HB2	1.94	0.50
1:F:338:MET:HG2	1:F:349:ILE:HG12	1.94	0.50
1:E:360:VAL:HG12	2:K:73:LEU:HD21	1.94	0.50
1:B:2:LYS:HB3	1:B:64:GLU:HB3	1.94	0.49
1:F:184:ILE:HD11	1:F:188:LEU:HD11	1.95	0.49
1:E:13:PRO:HA	1:E:230:PHE:HE1	1.77	0.49
1:D:100:ARG:HG2	1:D:105:ARG:HG3	1.94	0.49
1:E:2:LYS:HB3	1:E:64:GLU:HB2	1.94	0.49
1:D:338:MET:HE3	1:D:347:VAL:HG21	1.95	0.49
1:E:150:ASP:OD2	1:E:152:ARG:HD3	2.12	0.49
1:A:150:ASP:OD1	1:A:152:ARG:HD3	2.12	0.49
1:A:144:PHE:CD2	1:A:326:ASP:HB3	2.48	0.48
1:C:338:MET:HE3	1:C:347:VAL:HG21	1.95	0.48
1:C:255:HIS:HD2	1:C:339:MET:HG2	1.78	0.48
1:D:70:VAL:HG11	1:D:97:MET:SD	2.53	0.48
1:B:247:VAL:HG13	2:H:73:LEU:HD22	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:MET:HE3	1:B:347:VAL:HG21	1.96	0.48
1:C:170:VAL:HG22	1:C:179:VAL:HG23	1.96	0.48
1:A:53:MET:HB3	1:A:230:PHE:CZ	2.49	0.48
1:F:156:ASN:O	1:F:197:ARG:HB2	2.14	0.48
1:D:257:GLU:HG2	1:D:337:ARG:HG2	1.96	0.48
1:A:127:GLU:CD	1:C:39:ASP:H	2.17	0.47
1:D:170:VAL:HG22	1:D:179:VAL:HG23	1.95	0.47
1:C:144:PHE:CD2	1:C:326:ASP:HB3	2.49	0.47
1:F:150:ASP:OD2	1:F:152:ARG:HD3	2.15	0.47
1:A:21:LEU:H	1:A:21:LEU:HD12	1.80	0.47
1:A:338:MET:HE3	1:A:347:VAL:HG21	1.97	0.47
1:F:147:ALA:HB2	1:F:173:ASP:HA	1.97	0.47
1:F:33:LEU:HG	1:F:72:ALA:HB2	1.96	0.47
1:C:70:VAL:HG11	1:C:97:MET:SD	2.55	0.47
1:F:170:VAL:HG22	1:F:179:VAL:HG23	1.98	0.46
1:B:360:VAL:HG12	2:H:73:LEU:HD21	1.96	0.46
1:C:17:VAL:HG23	1:C:33:LEU:HD11	1.97	0.46
1:A:127:GLU:HG3	1:A:217:GLN:HG2	1.98	0.46
1:D:247:VAL:HG13	2:J:73:LEU:HD22	1.98	0.46
1:A:146:MET:HE1	1:A:197:ARG:HA	1.98	0.46
1:D:146:MET:HE1	1:D:197:ARG:HA	1.98	0.46
1:F:2:LYS:HD2	1:F:91:GLN:HB2	1.97	0.46
1:A:338:MET:HG2	1:A:349:ILE:HG12	1.97	0.46
1:B:33:LEU:HG	1:B:72:ALA:HB2	1.98	0.46
1:D:144:PHE:CD2	1:D:326:ASP:HB3	2.50	0.46
1:A:147:ALA:HB1	1:A:150:ASP:HB2	1.98	0.46
1:E:155:LEU:HD22	1:E:172:THR:HG23	1.97	0.45
1:E:33:LEU:HG	1:E:72:ALA:HB2	1.98	0.45
1:C:20:PRO:HG3	1:C:202:GLU:OE1	2.16	0.45
1:D:281:VAL:HG12	1:D:294:ALA:HB2	1.98	0.45
1:D:17:VAL:HG12	1:D:53:MET:HG3	1.97	0.45
1:F:33:LEU:O	1:F:69:THR:HA	2.16	0.45
1:E:144:PHE:CD2	1:E:326:ASP:HB3	2.51	0.45
1:A:50:GLU:HA	1:A:235:LYS:HD2	1.99	0.44
1:B:147:ALA:O	1:B:197:ARG:NH1	2.46	0.44
1:C:150:ASP:OD2	1:C:152:ARG:HD3	2.18	0.44
1:C:206:MET:HE2	1:C:227:VAL:HG12	1.99	0.44
1:C:155:LEU:HD22	1:C:172:THR:HG23	1.99	0.44
1:D:249:PRO:HD2	1:D:348:GLN:NE2	2.32	0.44
1:E:147:ALA:HB2	1:E:173:ASP:HA	1.98	0.44
1:E:174:GLY:HA2	2:K:74:PHE:CZ	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ALA:HB2	1:B:173:ASP:HA	1.98	0.44
1:A:33:LEU:HG	1:A:72:ALA:HB2	2.00	0.44
1:F:174:GLY:HA2	2:L:74:PHE:CZ	2.53	0.43
1:A:255[A]:HIS:NE2	1:A:339:MET:HG2	2.33	0.43
1:D:2:LYS:HG2	1:D:91:GLN:CB	2.48	0.43
1:B:16:GLN:HE21	1:B:53:MET:CE	2.29	0.43
1:C:147:ALA:HB2	1:C:173:ASP:HA	2.00	0.43
1:D:155:LEU:HD22	1:D:172:THR:HG23	2.01	0.43
1:E:147:ALA:HB1	1:E:150:ASP:HB2	2.01	0.43
1:E:70:VAL:HG11	1:E:97:MET:SD	2.58	0.43
1:A:263:LEU:HD23	1:A:336:VAL:HG21	1.99	0.43
1:F:346:SER:HA	1:F:363:PRO:HD3	2.01	0.43
1:B:147:ALA:HB1	1:B:150:ASP:HB2	2.00	0.43
1:B:70:VAL:HG11	1:B:97:MET:SD	2.58	0.43
1:D:27:LEU:HB2	1:D:30:LEU:HG	2.01	0.43
1:F:144:PHE:CD2	1:F:326:ASP:HB3	2.54	0.43
1:A:331:LEU:HD13	1:A:336:VAL:HG12	2.00	0.42
1:B:27:LEU:HB2	1:B:30:LEU:HG	2.01	0.42
1:B:338:MET:HG2	1:B:349:ILE:HG12	2.01	0.42
1:D:161:GLU:OE1	1:D:168:ARG:NH2	2.53	0.42
1:D:33:LEU:O	1:D:69:THR:HA	2.19	0.42
1:D:337:ARG:HG3	1:D:352:ALA:HA	2.02	0.42
1:A:14:LEU:O	1:A:18:SER:HB3	2.20	0.42
1:A:346:SER:HA	1:A:363:PRO:HD3	2.02	0.42
1:B:15:GLN:HG3	1:B:76:PHE:CE1	2.55	0.41
1:C:27:LEU:HB2	1:C:30:LEU:HG	2.01	0.41
1:C:12:LYS:HB3	1:C:13:PRO:HD3	2.01	0.41
1:A:159:LEU:O	1:A:169:THR:HA	2.21	0.41
1:A:27:LEU:HB2	1:A:30:LEU:HG	2.01	0.41
1:D:217:GLN:NE2	1:D:226:HIS:NE2	2.68	0.41
1:A:295:ASN:HA	1:A:300:GLU:O	2.21	0.41
1:D:156:ASN:O	1:D:197:ARG:HB2	2.21	0.41
1:D:285:VAL:HG12	1:D:290:LEU:HD13	2.03	0.41
1:A:283:LEU:O	1:A:316:GLU:HA	2.21	0.41
1:F:27:LEU:HB2	1:F:30:LEU:HG	2.03	0.41
1:E:159:LEU:O	1:E:169:THR:HA	2.22	0.40
1:F:339:MET:CE	1:F:350:GLU:HG2	2.51	0.40
1:E:249:PRO:HD2	1:E:348:GLN:NE2	2.30	0.40
1:A:156:ASN:O	1:A:197:ARG:HB2	2.21	0.40
1:B:50:GLU:HA	1:B:235:LYS:HD2	2.03	0.40
1:D:128:PHE:HB3	1:D:188:LEU:HD21	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:295:ASN:HA	1:F:300:GLU:O	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	365/366 (100%)	353 (97%)	11 (3%)	1 (0%)	41	50
1	B	360/366 (98%)	350 (97%)	9 (2%)	1 (0%)	41	50
1	C	363/366 (99%)	351 (97%)	12 (3%)	0	100	100
1	D	363/366 (99%)	355 (98%)	8 (2%)	0	100	100
1	E	360/366 (98%)	351 (98%)	8 (2%)	1 (0%)	41	50
1	F	363/366 (99%)	356 (98%)	7 (2%)	0	100	100
2	H	4/6 (67%)	3 (75%)	1 (25%)	0	100	100
2	I	4/6 (67%)	4 (100%)	0	0	100	100
2	J	4/6 (67%)	4 (100%)	0	0	100	100
2	K	4/6 (67%)	4 (100%)	0	0	100	100
2	L	4/6 (67%)	4 (100%)	0	0	100	100
All	All	2194/2226 (99%)	2135 (97%)	56 (3%)	3 (0%)	51	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	39	ASP
1	B	49	LEU
1	A	22	GLY

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	296/313 (95%)	291 (98%)	5 (2%)	60	76
1	B	303/313 (97%)	292 (96%)	11 (4%)	35	49
1	C	297/313 (95%)	287 (97%)	10 (3%)	37	51
1	D	297/313 (95%)	288 (97%)	9 (3%)	41	57
1	E	297/313 (95%)	291 (98%)	6 (2%)	55	72
1	F	303/313 (97%)	294 (97%)	9 (3%)	41	57
2	H	5/5 (100%)	4 (80%)	1 (20%)	1	1
2	I	5/5 (100%)	4 (80%)	1 (20%)	1	1
2	J	5/5 (100%)	4 (80%)	1 (20%)	1	1
2	K	5/5 (100%)	5 (100%)	0	100	100
2	L	5/5 (100%)	4 (80%)	1 (20%)	1	1
All	All	1818/1903 (96%)	1764 (97%)	54 (3%)	41	57

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

Mol	Chain	Res	Type
1	A	129	THR
1	A	176	ARG
1	A	188	LEU
1	A	192	SER
1	A	212	ASN
1	B	26	THR
1	B	39	ASP
1	B	56	ARG
1	B	125	GLU
1	B	129	THR
1	B	176	ARG
1	B	184	ILE
1	B	191	HIS
1	B	212	ASN
1	B	253	ASP

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Mol	Chain	Res	Type
1	B	295	ASN
1	C	21	LEU
1	C	39	ASP
1	C	87	GLU
1	C	129	THR
1	C	176	ARG
1	C	192	SER
1	C	212	ASN
1	C	253	ASP
1	C	285	VAL
1	C	355	GLN
1	D	56	ARG
1	D	93	GLU
1	D	121	ASP
1	D	123	GLN
1	D	129	THR
1	D	176	ARG
1	D	253	ASP
1	D	334	GLU
1	D	364	MET
1	E	64	GLU
1	E	129	THR
1	E	176	ARG
1	E	192	SER
1	E	334	GLU
1	E	355	GLN
1	F	11	LEU
1	F	27	LEU
1	F	62	PRO
1	F	84	GLU
1	F	129	THR
1	F	176	ARG
1	F	192	SER
1	F	251	ASN
1	F	285	VAL
2	H	72	ASP
2	I	72	ASP
2	J	72	ASP
2	L	72	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	61	GLN
1	A	191	HIS
1	A	217	GLN
1	A	335	ASN
1	A	348	GLN
1	B	16	GLN
1	B	61	GLN
1	B	143	GLN
1	B	217	GLN
1	B	265	GLN
1	B	335	ASN
1	B	348	GLN
1	C	9	HIS
1	C	91	GLN
1	C	191	HIS
1	C	255	HIS
1	C	265	GLN
1	C	335	ASN
1	C	348	GLN
1	D	16	GLN
1	D	186	GLN
1	D	191	HIS
1	D	217	GLN
1	D	265	GLN
1	D	335	ASN
1	D	348	GLN
1	E	191	HIS
1	E	217	GLN
1	E	265	GLN
1	E	335	ASN
1	E	348	GLN
1	F	9	HIS
1	F	191	HIS
1	F	217	GLN
1	F	265	GLN
1	F	335	ASN
1	F	348	GLN
2	I	70	GLN
2	J	70	GLN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	366/366 (100%)	0.31	20 (5%) 25 31	27, 57, 85, 104	1 (0%)
1	B	364/366 (99%)	0.00	7 (1%) 66 73	23, 45, 72, 97	2 (0%)
1	C	365/366 (99%)	0.24	21 (5%) 23 29	31, 55, 80, 101	2 (0%)
1	D	365/366 (99%)	0.23	11 (3%) 50 57	32, 54, 83, 95	1 (0%)
1	E	364/366 (99%)	0.11	13 (3%) 42 49	26, 51, 76, 107	1 (0%)
1	F	365/366 (99%)	-0.04	9 (2%) 57 64	29, 47, 75, 99	1 (0%)
2	H	5/6 (83%)	0.42	0 100 100	44, 45, 58, 59	0
2	I	5/6 (83%)	2.05	3 (60%) 0 0	62, 67, 74, 78	0
2	J	5/6 (83%)	2.99	3 (60%) 0 0	78, 90, 100, 108	0
2	K	5/6 (83%)	0.90	0 100 100	54, 54, 65, 66	0
2	L	5/6 (83%)	1.94	3 (60%) 0 0	66, 72, 81, 86	0
All	All	2214/2226 (99%)	0.16	90 (4%) 37 44	23, 52, 81, 108	8 (0%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	23	GLY	7.4
1	D	23	GLY	7.1
1	A	22	GLY	6.8
1	A	366	LEU	5.4
2	J	74	PHE	5.4
1	A	23	GLY	5.3
1	C	22	GLY	5.2
1	A	21	LEU	5.2
1	A	153	TYR	5.1
1	D	24	ARG	5.0
1	C	60	VAL	4.9
1	D	211	ASP	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	187	SER	4.1
1	E	24	ARG	4.1
1	F	26	THR	4.0
1	B	20	PRO	4.0
1	D	22	GLY	3.9
2	L	74	PHE	3.8
1	E	94	GLY	3.8
2	I	71	LEU	3.7
2	J	71	LEU	3.7
1	E	211	ASP	3.5
1	C	40	GLY	3.4
1	A	20	PRO	3.4
1	E	121	ASP	3.4
2	J	72	ASP	3.4
1	B	24	ARG	3.3
1	B	26	THR	3.3
1	E	208	ASP	3.2
1	D	315	MET	3.1
1	E	153	TYR	3.1
2	I	74	PHE	3.1
1	E	210	GLY	3.0
1	F	23	GLY	3.0
1	C	114	ALA	3.0
2	L	72	ASP	3.0
1	C	255	HIS	3.0
1	F	22	GLY	2.9
1	A	148	HIS	2.9
1	C	85	GLY	2.9
1	A	311	SER	2.8
1	C	58	ALA	2.7
1	E	114	ALA	2.7
1	F	24	ARG	2.7
1	A	246	ARG	2.7
1	F	299	GLN	2.7
1	B	187	SER	2.7
1	D	209	GLY	2.7
1	A	133	ALA	2.6
1	D	120	ASP	2.6
1	D	208	ASP	2.6
1	F	118	ASN	2.6
1	C	308	VAL	2.5
1	E	21	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	311	SER	2.5
1	F	149	GLN	2.5
1	A	120	ASP	2.5
1	E	209	GLY	2.4
1	B	60	VAL	2.4
1	C	26	THR	2.4
1	A	209	GLY	2.4
1	D	25	PRO	2.4
1	B	84	GLU	2.4
1	C	117	PRO	2.3
1	A	60	VAL	2.3
1	A	26	THR	2.3
1	E	187	SER	2.3
1	C	119	LEU	2.3
1	E	26	THR	2.3
1	C	120	ASP	2.3
1	A	149	GLN	2.3
1	B	238	ASP	2.2
1	E	39	ASP	2.2
2	L	71	LEU	2.2
1	D	362	MET	2.2
1	C	313	ALA	2.2
1	C	35	LEU	2.2
1	C	62	PRO	2.2
1	D	148	HIS	2.2
1	F	251	ASN	2.1
1	A	191	HIS	2.1
1	C	151	VAL	2.1
1	C	353	ALA	2.1
2	I	72	ASP	2.1
1	A	365	ARG	2.1
1	F	250	LYS	2.1
1	A	24	ARG	2.1
1	A	208	ASP	2.1
1	C	55	ALA	2.0
1	C	310	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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.