



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:54 PM UTC

PDB ID : 6DVD / pdb_00006dvd
Title : Crystal structure of Mycobacterium tuberculosis transcription initiation complex(ECF sigma factor L) with 6 nt spacer and bromine labelled in position "-11
Authors : Lin, W.; Das, K.; Feng, Y.; Ebright, R.H.
Deposited on : 2018-06-23
Resolution : 3.90 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 : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

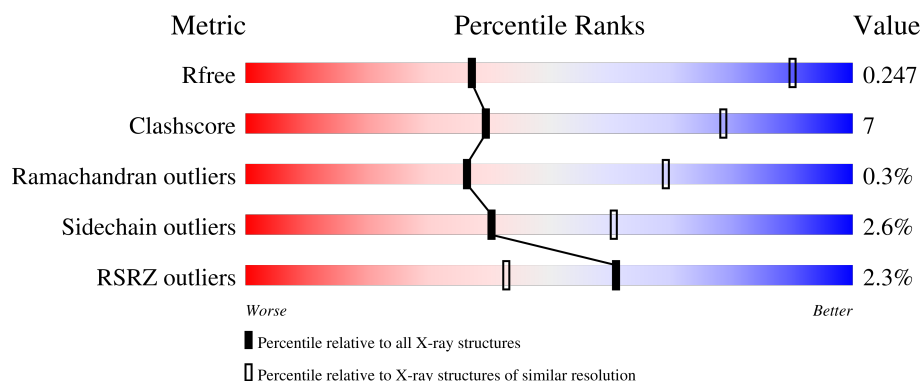
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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}	180053	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 $\leq 5\%$. 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	359	50% 12% • 37%
1	B	359	2% 48% 15% • 35%
2	C	1178	2% 79% 16% • •
3	D	1316	• 79% 16% •
4	E	110	5% 64% 10% 26%

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Mol	Chain	Length	Quality of chain
5	F	177	7% 84% 14% ••
6	G	10	50% 50%
7	H	24	25% 38% 63%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 24744 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-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1716	1080	296	338	2			
1	B	232	Total	C	N	O	S	0	0	0
			1732	1093	296	341	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP P9WGZ1
A	-10	GLY	-	expression tag	UNP P9WGZ1
A	-9	HIS	-	expression tag	UNP P9WGZ1
A	-8	HIS	-	expression tag	UNP P9WGZ1
A	-7	HIS	-	expression tag	UNP P9WGZ1
A	-6	HIS	-	expression tag	UNP P9WGZ1
A	-5	HIS	-	expression tag	UNP P9WGZ1
A	-4	HIS	-	expression tag	UNP P9WGZ1
A	-3	HIS	-	expression tag	UNP P9WGZ1
A	-2	HIS	-	expression tag	UNP P9WGZ1
A	-1	HIS	-	expression tag	UNP P9WGZ1
A	0	HIS	-	expression tag	UNP P9WGZ1
B	-11	MET	-	initiating methionine	UNP P9WGZ1
B	-10	GLY	-	expression tag	UNP P9WGZ1
B	-9	HIS	-	expression tag	UNP P9WGZ1
B	-8	HIS	-	expression tag	UNP P9WGZ1
B	-7	HIS	-	expression tag	UNP P9WGZ1
B	-6	HIS	-	expression tag	UNP P9WGZ1
B	-5	HIS	-	expression tag	UNP P9WGZ1
B	-4	HIS	-	expression tag	UNP P9WGZ1
B	-3	HIS	-	expression tag	UNP P9WGZ1
B	-2	HIS	-	expression tag	UNP P9WGZ1
B	-1	HIS	-	expression tag	UNP P9WGZ1
B	0	HIS	-	expression tag	UNP P9WGZ1

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1126	Total	C	N	O	S	0	0	0
			8724	5459	1531	1695	39			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1265	Total	C	N	O	S	0	0	0
			9895	6195	1794	1866	40			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	81	Total	C	N	O	0	0	0
			630	403	106	121			

- Molecule 5 is a protein called ECF RNA polymerase sigma factor SigL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	174	Total	C	N	O	S	0	0	0
			1352	840	256	254	2			

- Molecule 6 is a DNA chain called DNA (5'-D(*GP*CP*AP*TP*CP*CP*GP*TP*GP*A)-3').

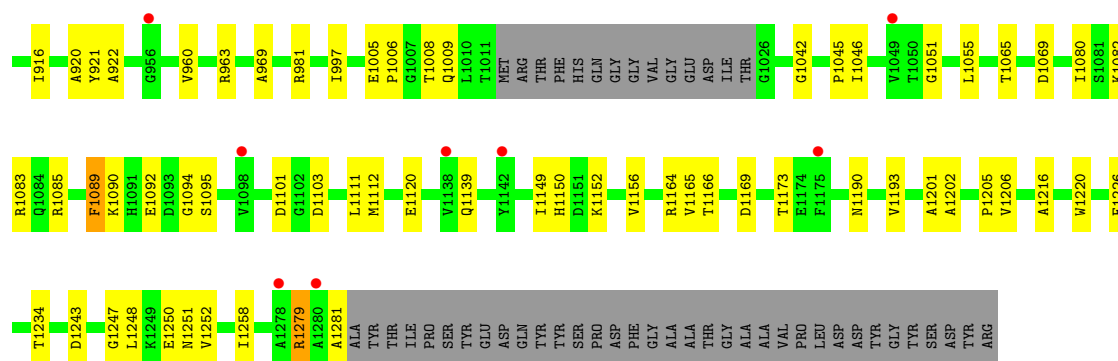
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	10	Total	C	N	O	P	0	0	0
			202	97	38	58	9			

- Molecule 7 is a DNA chain called DNA (5'-D(P*CP*GP*TP*GP*TP*CP*AP*GP*TP*AP*GP*CP*TP*GP*TP*CP*AP*CP*GP*GP*AP*TP*GP*C)-3').

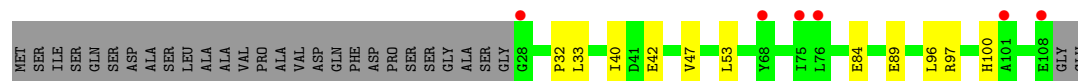
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	24	Total	C	N	O	P	0	0	0
			491	234	90	144	23			

- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

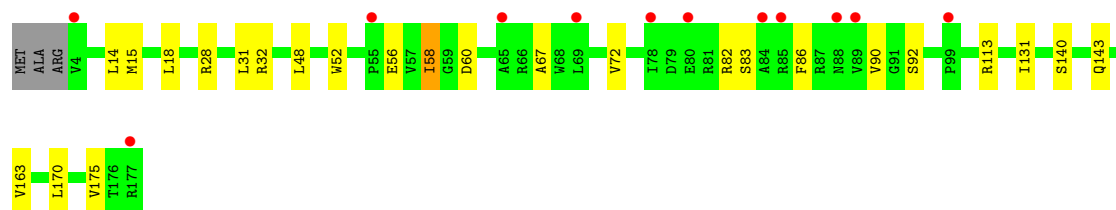
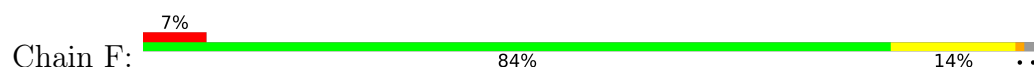
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	D	2	Total	Zn	0	0
			2	2		



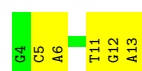
- Molecule 4: DNA-directed RNA polymerase subunit omega



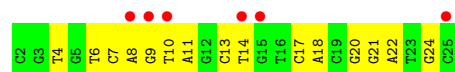
- Molecule 5: ECF RNA polymerase sigma factor SigL



- Molecule 6: DNA (5'-D(*GP*CP*AP*TP*CP*CP*GP*TP*GP*A)-3')



- Molecule 7: DNA (5'-D(P*CP*GP*TP*GP*TP*CP*AP*GP*TP*AP*GP*CP*TP*GP*TP*C P*AP*CP*GP*GP*AP*TP*GP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	142.13Å 161.49Å 239.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.90 – 3.90 45.90 – 3.90	Depositor EDS
% Data completeness (in resolution range)	73.8 (45.90-3.90) 73.8 (45.90-3.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 3.88Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.220 , 0.246 0.222 , 0.247	Depositor DCC
R_{free} test set	2000 reflections (5.31%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.956	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 56.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	24744	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.51% 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: ZN

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	0.09	0/1742	0.31	0/2370
1	B	0.10	0/1758	0.31	0/2397
2	C	0.09	0/8883	0.30	0/12043
3	D	0.09	0/10061	0.28	0/13600
4	E	0.09	0/643	0.26	0/877
5	F	0.11	0/1374	0.30	0/1869
6	G	0.16	0/226	0.33	0/347
7	H	0.21	0/550	0.50	0/848
All	All	0.10	0/25237	0.30	0/34351

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 [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	1716	0	1756	30	0
1	B	1732	0	1754	37	0
2	C	8724	0	8651	121	0
3	D	9895	0	9954	144	0
4	E	630	0	622	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	F	1352	0	1346	19	0
6	G	202	0	114	7	0
7	H	491	0	272	19	0
8	D	2	0	0	0	0
All	All	24744	0	24469	327	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 (327) 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:40:ARG:HE	1:B:33:THR:HG22	1.46	0.81
3:D:334:ARG:HD3	5:F:90:VAL:HG21	1.63	0.80
1:A:4:SER:HB3	1:B:144:ARG:HH12	1.48	0.76
3:D:363:PRO:HG2	5:F:15:MET:HG3	1.67	0.75
3:D:454:PRO:HA	3:D:457:MET:HE2	1.70	0.74
3:D:1090:LYS:HE2	3:D:1103:ASP:HA	1.69	0.74
2:C:593:MET:HA	2:C:628:THR:HG21	1.70	0.73
3:D:832:ILE:HG22	3:D:834:ARG:H	1.54	0.72
2:C:541:VAL:HG12	2:C:578:TYR:HB2	1.71	0.72
2:C:1024:THR:H	3:D:730:THR:HG21	1.52	0.72
1:B:84:VAL:HG12	1:B:199:LYS:HD2	1.72	0.71
3:D:366:ILE:HD11	5:F:15:MET:HE2	1.72	0.71
1:B:90:ASP:OD1	1:B:90:ASP:N	2.18	0.71
2:C:558:ARG:HB3	2:C:570:TYR:HB3	1.72	0.70
5:F:14:LEU:HD21	5:F:58:ILE:HG21	1.72	0.70
6:G:12:DG:H2''	6:G:13:DA:H5'	1.74	0.70
3:D:1247:GLY:O	3:D:1251:ASN:ND2	2.23	0.70
2:C:454:ARG:HH12	2:C:487:GLU:HG2	1.58	0.69
3:D:1090:LYS:HB3	3:D:1092:GLU:HG2	1.74	0.68
3:D:1055:LEU:H	3:D:1101:ASP:HB3	1.59	0.67
2:C:928:ILE:HD11	3:D:817:LEU:HD11	1.76	0.66
1:B:81:LYS:HD3	1:B:165:ASP:HB2	1.79	0.65
3:D:1045:PRO:HG2	3:D:1111:LEU:HB2	1.79	0.65
3:D:1248:LEU:HD22	3:D:1258:ILE:HB	1.79	0.65
3:D:882:GLN:HG3	3:D:997:ILE:HD11	1.79	0.64
2:C:926:MET:HE1	3:D:817:LEU:HA	1.79	0.64
2:C:536:GLU:OE2	2:C:562:ARG:NH2	2.31	0.64
2:C:104:SER:HB3	2:C:140:ILE:HB	1.80	0.63
3:D:1051:GLY:HA2	3:D:1069:ASP:HB2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:122:CYS:HA	2:C:127:MET:HG3	1.79	0.63
3:D:458:LYS:NZ	3:D:462:ASP:OD2	2.31	0.63
3:D:50:LYS:HE2	3:D:79:GLY:HA3	1.82	0.62
6:G:11:DT:H2'	6:G:12:DG:C8	2.35	0.62
1:B:100:GLN:HG3	1:B:133:LYS:HB2	1.82	0.61
3:D:104:ILE:HD12	3:D:379:ASP:HB3	1.83	0.61
2:C:211:TRP:HZ3	7:H:14:DT:H3'	1.65	0.61
2:C:401:ARG:HA	2:C:404:MET:HE2	1.81	0.61
1:A:186:ARG:HG3	1:A:187:THR:HG23	1.84	0.60
2:C:168:ILE:HB	2:C:173:ARG:HD2	1.83	0.60
1:A:83:LEU:HA	1:A:123:MET:HE1	1.84	0.60
3:D:436:LEU:HD11	3:D:523:GLN:HB3	1.84	0.59
1:A:40:ARG:NH2	2:C:1013:GLY:O	2.33	0.59
2:C:234:VAL:HG12	2:C:261:THR:HG21	1.84	0.59
1:B:55:ARG:NH2	1:B:137:GLU:OE1	2.36	0.59
3:D:1173:THR:HG22	3:D:1193:VAL:HG21	1.84	0.59
2:C:524:VAL:HG21	2:C:548:ILE:HD13	1.84	0.58
1:A:144:ARG:NH2	1:B:27:GLU:OE2	2.36	0.58
1:B:92:PRO:HB3	1:B:141:GLU:HG2	1.86	0.58
1:B:24:GLU:HB2	1:B:191:LYS:HG3	1.85	0.58
2:C:209:GLY:O	7:H:13:DC:N4	2.37	0.58
2:C:571:VAL:HG22	2:C:572:PRO:HD2	1.86	0.58
3:D:1092:GLU:HG3	3:D:1094:GLY:H	1.68	0.58
2:C:824:ILE:HA	5:F:163:VAL:HG13	1.86	0.58
3:D:600:GLN:HB2	3:D:609:THR:HB	1.85	0.57
2:C:32:VAL:HG13	2:C:33:PRO:HD3	1.86	0.57
3:D:337:THR:OG1	3:D:341:ASN:ND2	2.33	0.57
3:D:1165:VAL:HG12	3:D:1205:PRO:HA	1.87	0.57
3:D:190:LYS:HE3	3:D:192:ASP:HB3	1.86	0.57
3:D:1045:PRO:HB2	3:D:1111:LEU:HD12	1.87	0.57
5:F:31:LEU:HD21	7:H:10:DT:C2	2.40	0.57
1:B:147:VAL:HG13	1:B:166:SER:HB2	1.87	0.57
2:C:113:ASP:HB3	2:C:132:PRO:HG2	1.87	0.57
2:C:47:PRO:HG2	2:C:581:VAL:HG13	1.88	0.56
2:C:1087:GLU:HG3	2:C:1091:ILE:HD11	1.85	0.56
1:A:216:VAL:HG13	1:B:216:VAL:HG13	1.88	0.56
2:C:607:MET:SD	2:C:892:LYS:NZ	2.78	0.56
3:D:107:PHE:HZ	3:D:126:GLU:HG2	1.71	0.56
2:C:83:VAL:HG13	2:C:87:GLU:HB2	1.87	0.55
2:C:398:ARG:HH12	5:F:32:ARG:HB3	1.71	0.55
3:D:789:LEU:HD22	3:D:793:TYR:HE2	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:826:ASN:HB3	3:D:832:ILE:HD11	1.88	0.55
1:A:22:VAL:HG12	1:A:193:ILE:HG12	1.88	0.55
2:C:458:LEU:HD21	2:C:496:LEU:HD13	1.88	0.55
2:C:1055:GLN:HG2	2:C:1094:ASP:HB3	1.88	0.55
3:D:1279:ARG:HE	3:D:1281:ALA:H	1.54	0.55
3:D:1083:ARG:HD2	3:D:1112:MET:HE3	1.88	0.55
3:D:890:ASP:OD1	3:D:963:ARG:NH2	2.33	0.55
3:D:573:PRO:HB2	3:D:576:MET:HE2	1.89	0.54
1:B:17:ASN:OD1	1:B:17:ASN:N	2.39	0.54
2:C:982:GLU:HG3	3:D:841:ARG:HH12	1.72	0.54
2:C:642:VAL:HB	2:C:703:ALA:HB3	1.89	0.54
3:D:1042:GLY:O	3:D:1083:ARG:NH2	2.37	0.54
2:C:853:PHE:HD2	2:C:868:LEU:HD23	1.72	0.53
2:C:731:TYR:HE1	3:D:579:LEU:HB2	1.73	0.53
2:C:298:ASN:HA	2:C:302:LYS:HB2	1.91	0.53
2:C:369:ASP:C	2:C:371:ASP:H	2.16	0.53
2:C:168:ILE:HG12	2:C:431:PHE:HB3	1.91	0.53
3:D:443:LEU:HD13	3:D:448:ALA:HB2	1.90	0.53
3:D:729:VAL:HG11	3:D:802:ILE:HD11	1.91	0.53
3:D:1164:ARG:NH2	3:D:1216:ALA:O	2.42	0.53
3:D:63:GLY:HA2	3:D:66:LYS:HE2	1.91	0.53
2:C:217:ASP:OD2	2:C:231:ARG:NH2	2.42	0.52
3:D:585:LEU:O	3:D:589:THR:OG1	2.27	0.52
1:B:129:ASN:OD1	1:B:130:ASP:N	2.38	0.52
3:D:589:THR:HG21	3:D:688:MET:HG2	1.91	0.52
3:D:34:ILE:HA	3:D:41:PRO:HA	1.91	0.52
5:F:67:ALA:O	7:H:8:DA:N6	2.43	0.52
2:C:960:PRO:HD2	2:C:963:LEU:HD22	1.91	0.52
2:C:190:THR:HG23	2:C:199:LEU:HB2	1.91	0.51
2:C:704:ASP:HB2	2:C:708:THR:HB	1.91	0.51
3:D:117:LEU:HD12	3:D:299:VAL:HG22	1.92	0.51
2:C:257:ILE:HD11	2:C:346:VAL:HG23	1.92	0.51
6:G:12:DG:C2'	6:G:13:DA:H5'	2.40	0.51
1:A:30:PHE:HE1	1:B:41:THR:HA	1.74	0.51
1:B:102:PRO:HD3	1:B:131:LYS:H	1.75	0.51
1:B:148:PRO:HG3	3:D:626:VAL:HG21	1.93	0.51
2:C:211:TRP:CZ3	7:H:14:DT:H3'	2.45	0.51
2:C:1045:SER:HB3	3:D:450:GLU:O	2.10	0.51
1:B:146:TYR:O	3:D:624:ARG:NE	2.44	0.51
2:C:853:PHE:HB2	2:C:868:LEU:HB3	1.93	0.51
1:A:56:ILE:HB	1:A:59:VAL:HG22	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:VAL:HG13	1:A:119:HIS:HB2	1.93	0.51
2:C:740:ARG:NH1	2:C:744:GLU:OE1	2.44	0.51
2:C:451:HIS:HA	2:C:454:ARG:HG2	1.91	0.51
3:D:505:HIS:CD2	3:D:507:LEU:HB2	2.45	0.51
1:A:7:PRO:HA	1:A:25:PRO:HD2	1.93	0.50
2:C:377:ARG:NH2	2:C:383:GLU:OE1	2.40	0.50
2:C:182:SER:HB2	2:C:377:ARG:HB2	1.94	0.50
2:C:446:LEU:HB2	2:C:713:MET:HE2	1.93	0.50
3:D:447:MET:HE1	3:D:543:VAL:HG21	1.94	0.50
3:D:921:TYR:O	3:D:1150:HIS:NE2	2.41	0.50
5:F:72:VAL:HG22	7:H:4:DT:C2	2.47	0.50
4:E:89:GLU:OE2	4:E:97:ARG:NH1	2.45	0.50
2:C:686:GLN:HA	2:C:705:GLY:HA2	1.94	0.49
3:D:1080:ILE:HG22	3:D:1082:LYS:H	1.76	0.49
3:D:173:ARG:HH21	3:D:201:GLY:HA2	1.78	0.49
6:G:5:DC:H2"	6:G:6:DA:C8	2.47	0.49
1:A:98:ARG:HG3	1:A:135:GLU:HG3	1.95	0.49
1:B:102:PRO:HB3	1:B:130:ASP:HA	1.94	0.49
3:D:460:LEU:HD11	3:D:483:VAL:HG12	1.94	0.49
3:D:778:TRP:CD2	3:D:835:PRO:HG3	2.47	0.49
1:A:89:GLU:HB3	1:A:91:GLU:HG2	1.95	0.49
3:D:500:ARG:HD2	3:D:534:ALA:HB2	1.95	0.49
5:F:28:ARG:HB3	7:H:9:DG:H2"	1.94	0.49
5:F:140:SER:HB3	5:F:143:GLN:HG3	1.94	0.49
2:C:751:HIS:CD2	2:C:877:ARG:HD2	2.48	0.49
2:C:1052:ILE:O	3:D:89:ARG:NH1	2.45	0.49
1:A:217:GLU:HG2	1:B:232:ILE:HD12	1.95	0.48
2:C:1067:ARG:HA	3:D:421:ARG:HA	1.95	0.48
7:H:6:DT:H3'	7:H:7:DC:H5"	1.95	0.48
2:C:62:GLU:OE1	2:C:69:ARG:NE	2.45	0.48
2:C:446:LEU:HD13	2:C:593:MET:HE1	1.94	0.48
3:D:1220:TRP:CD1	3:D:1243:ASP:HB2	2.48	0.48
4:E:42:GLU:OE1	4:E:100:HIS:NE2	2.43	0.48
1:B:98:ARG:HG2	1:B:135:GLU:HG2	1.96	0.48
2:C:322:LEU:HD23	2:C:357:VAL:HG21	1.95	0.48
3:D:557:ILE:HG23	4:E:40:ILE:HD11	1.96	0.48
2:C:473:ARG:NH2	2:C:492:PRO:O	2.46	0.48
3:D:1139:GLN:NE2	3:D:1149:ILE:O	2.40	0.48
5:F:18:LEU:HD13	5:F:48:LEU:HD22	1.95	0.48
2:C:450:THR:HG21	2:C:613:ARG:HD2	1.95	0.48
3:D:1247:GLY:H	3:D:1251:ASN:HD21	1.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:189:GLU:HB2	2:C:367:THR:HG21	1.96	0.48
2:C:239:LYS:NZ	2:C:265:ASP:OD2	2.47	0.48
2:C:730:ASN:HA	2:C:734:ALA:HB3	1.94	0.48
2:C:848:ILE:HD13	2:C:874:ALA:HB2	1.95	0.48
3:D:505:HIS:ND1	3:D:1005:GLU:HG3	2.29	0.47
2:C:369:ASP:O	2:C:370:ILE:HG12	2.14	0.47
2:C:388:GLN:HG3	2:C:430:PHE:HB2	1.96	0.47
3:D:1169:ASP:H	3:D:1202:ALA:HB3	1.79	0.47
2:C:549:ASP:OD1	2:C:550:ALA:N	2.47	0.47
2:C:974:THR:HG23	2:C:980:ALA:H	1.79	0.47
1:A:188:ASP:OD2	1:B:151:GLN:NE2	2.44	0.47
1:B:170:PRO:HB2	1:B:202:ILE:HD11	1.97	0.47
2:C:795:GLU:HG2	2:C:846:LYS:HG2	1.96	0.47
3:D:473:LYS:NZ	3:D:477:GLU:OE2	2.45	0.47
4:E:47:VAL:HG11	4:E:53:LEU:HB2	1.97	0.47
3:D:500:ARG:HB2	3:D:541:MET:HG2	1.95	0.47
3:D:882:GLN:HG2	3:D:1248:LEU:HB2	1.95	0.47
3:D:384:ASN:H	3:D:401:SER:HB3	1.80	0.47
3:D:611:VAL:HG22	3:D:634:LYS:HB2	1.97	0.46
3:D:872:TYR:CD1	6:G:13:DA:H5''	2.51	0.46
2:C:731:TYR:CE1	3:D:579:LEU:HB2	2.49	0.46
7:H:10:DT:H2'	7:H:11:DA:C8	2.50	0.46
2:C:602:ALA:HB2	3:D:853:THR:HG22	1.96	0.46
2:C:1119:GLU:OE2	3:D:89:ARG:NH2	2.48	0.46
2:C:1146:GLU:H	2:C:1146:GLU:HG2	1.48	0.46
3:D:1247:GLY:H	3:D:1251:ASN:ND2	2.12	0.46
1:A:56:ILE:HG12	1:A:136:VAL:HG22	1.98	0.46
3:D:340:LEU:HD11	3:D:405:LEU:HD11	1.98	0.46
3:D:872:TYR:CG	6:G:13:DA:H5''	2.50	0.46
3:D:922:ALA:HB2	3:D:981:ARG:HH21	1.81	0.46
6:G:12:DG:H2'	6:G:13:DA:C8	2.50	0.46
7:H:9:DG:N3	7:H:9:DG:H2'	2.31	0.46
3:D:431:VAL:O	3:D:523:GLN:HA	2.15	0.46
3:D:1085:ARG:HA	3:D:1112:MET:HA	1.98	0.46
7:H:7:DC:OP1	7:H:7:DC:H4'	2.15	0.46
2:C:628:THR:HG23	2:C:630:MET:H	1.80	0.46
2:C:38:ARG:HA	2:C:971:ILE:HG13	1.98	0.46
2:C:588:SER:OG	2:C:589:VAL:N	2.48	0.46
3:D:105:TRP:HB3	3:D:1234:THR:HG22	1.98	0.46
1:A:18:ARG:NH2	2:C:997:ASP:OD1	2.48	0.45
2:C:486:ILE:HD11	3:D:849:TYR:HE2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:412:ILE:HD13	2:C:417:LEU:HD21	1.98	0.45
3:D:1166:THR:HB	3:D:1206:VAL:HG21	1.98	0.45
3:D:585:LEU:HD13	3:D:673:PHE:HE2	1.81	0.45
2:C:32:VAL:H	2:C:33:PRO:CD	2.29	0.45
3:D:235:ILE:HD12	3:D:241:TYR:HD1	1.81	0.45
2:C:32:VAL:H	2:C:33:PRO:HD3	1.81	0.45
2:C:115:VAL:HG11	2:C:129:TYR:CZ	2.52	0.45
2:C:522:GLY:O	2:C:553:ARG:HA	2.16	0.45
2:C:901:VAL:HG23	2:C:912:PRO:HG3	1.98	0.45
1:A:40:ARG:NE	1:B:33:THR:HG22	2.25	0.45
2:C:544:ALA:HB2	2:C:580:ASP:HB2	1.99	0.44
3:D:59:GLU:HG2	3:D:66:LYS:HD3	1.98	0.44
3:D:463:LEU:HB2	3:D:465:HIS:HD2	1.82	0.44
3:D:589:THR:HG22	3:D:670:ARG:HG2	1.99	0.44
3:D:1190:ASN:ND2	3:D:1201:ALA:O	2.44	0.44
2:C:313:ARG:HH22	2:C:337:ASP:CG	2.25	0.44
1:B:83:LEU:HB2	1:B:123:MET:HE1	1.98	0.44
2:C:213:GLU:OE1	2:C:225:ARG:NH1	2.50	0.44
2:C:1089:LEU:HD13	3:D:1252:VAL:HG13	1.99	0.44
3:D:453:LYS:O	3:D:457:MET:HG3	2.17	0.44
3:D:468:ASN:HD22	5:F:175:VAL:HG22	1.82	0.44
3:D:575:ALA:O	3:D:713:VAL:HG21	2.18	0.44
1:B:84:VAL:HG23	1:B:119:HIS:HB2	2.00	0.44
3:D:103:HIS:HB3	3:D:106:TYR:HD2	1.82	0.44
1:A:95:MET:HE3	1:A:110:ILE:HG21	2.00	0.44
2:C:927:ASN:O	2:C:930:GLN:HG2	2.17	0.44
2:C:680:HIS:O	2:C:1034:HIS:HE1	2.01	0.44
2:C:809:LYS:HE2	2:C:813:GLU:HB2	2.00	0.44
2:C:1104:GLU:OE1	5:F:113:ARG:NH1	2.50	0.44
3:D:588:LEU:HD12	3:D:589:THR:HG23	2.00	0.44
1:A:173:LYS:HE3	1:A:173:LYS:HB2	1.84	0.43
2:C:484:CYS:HB2	2:C:588:SER:HB2	2.00	0.43
3:D:337:THR:O	5:F:92:SER:HA	2.18	0.43
2:C:454:ARG:HH11	2:C:499:SER:HB2	1.82	0.43
3:D:291:ARG:NH2	7:H:24:DG:O6	2.44	0.43
2:C:140:ILE:HA	2:C:147:ILE:HG12	1.99	0.43
2:C:228:ARG:O	2:C:228:ARG:HG3	2.17	0.43
2:C:344:TYR:OH	2:C:364:PRO:O	2.25	0.43
2:C:513:GLU:HB3	2:C:530:TYR:HB3	2.01	0.43
2:C:737:LEU:HG	2:C:895:ILE:HD12	1.99	0.43
2:C:982:GLU:HG3	3:D:841:ARG:NH1	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:33:THR:HB	3:D:327:MET:HE1	2.00	0.43
1:B:24:GLU:HA	1:B:25:PRO:HA	1.82	0.43
3:D:505:HIS:HD2	3:D:507:LEU:H	1.64	0.43
3:D:445:LYS:HB3	3:D:484:TRP:CZ3	2.53	0.43
1:A:81:LYS:NZ	1:A:165:ASP:HB2	2.34	0.43
1:B:74:THR:HG21	3:D:608:GLU:HB2	2.00	0.43
2:C:455:LEU:HD12	2:C:483:MET:HG3	2.00	0.43
3:D:579:LEU:HD23	3:D:579:LEU:HA	1.88	0.43
3:D:893:THR:HG21	3:D:969:ALA:HB3	2.00	0.43
1:B:28:PRO:HA	1:B:29:GLY:HA2	1.47	0.43
3:D:123:LYS:HE3	3:D:127:LYS:HE2	2.00	0.43
3:D:1279:ARG:H	3:D:1279:ARG:HG3	1.64	0.43
5:F:82:ARG:HB3	5:F:83:SER:H	1.67	0.43
3:D:453:LYS:HG3	3:D:476:VAL:HG11	2.00	0.43
2:C:86:LEU:HD23	2:C:86:LEU:HA	1.90	0.43
2:C:474:ASP:HA	3:D:857:ARG:HD2	2.00	0.43
3:D:24:SER:HB2	3:D:94:HIS:HB3	2.01	0.43
3:D:262:GLN:HB2	3:D:313:VAL:HG11	2.01	0.42
3:D:294:LYS:NZ	7:H:22:DA:O5'	2.52	0.42
3:D:859:GLY:O	3:D:863:THR:OG1	2.32	0.42
5:F:170:LEU:HD22	5:F:175:VAL:HG21	2.01	0.42
1:B:30:PHE:HA	1:B:33:THR:HG23	2.01	0.42
2:C:592:ALA:HA	2:C:976:VAL:HG21	2.01	0.42
2:C:731:TYR:HB3	3:D:432:VAL:HG21	2.00	0.42
2:C:906:PHE:HA	2:C:912:PRO:HA	2.01	0.42
3:D:16:THR:HG22	3:D:18:GLU:H	1.84	0.42
3:D:867:THR:HG22	3:D:1008:THR:HG23	2.01	0.42
4:E:96:LEU:HD12	4:E:96:LEU:HA	1.86	0.42
3:D:139:VAL:HA	3:D:252:PHE:HA	2.00	0.42
3:D:670:ARG:NH1	3:D:685:ASN:O	2.52	0.42
1:A:3:ILE:HG13	1:A:4:SER:H	1.85	0.42
2:C:57:GLN:HG3	2:C:452:LYS:HB3	2.01	0.42
3:D:128:ILE:HD11	3:D:234:LEU:HD11	2.01	0.42
3:D:427:ARG:NH1	3:D:540:GLN:OE1	2.52	0.42
2:C:610:ASN:OD1	2:C:613:ARG:NH1	2.52	0.42
3:D:598:GLU:HG2	3:D:599:TYR:H	1.84	0.42
3:D:20:ILE:HG23	3:D:318:PRO:HB3	2.02	0.42
3:D:97:LEU:HD22	3:D:374:LEU:HD21	2.00	0.42
3:D:577:PRO:HA	3:D:581:MET:HE3	2.02	0.42
1:A:206:ASP:OD1	1:B:226:ASN:ND2	2.52	0.42
2:C:56:VAL:HG21	2:C:500:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:120:ASP:OD1	2:C:120:ASP:N	2.53	0.42
2:C:228:ARG:HH21	7:H:13:DC:H1'	1.83	0.42
2:C:305:ARG:HA	2:C:305:ARG:HD2	1.83	0.42
2:C:767:GLU:HG2	2:C:807:THR:HG22	2.02	0.42
3:D:447:MET:HE2	3:D:447:MET:HB3	1.88	0.42
3:D:897:ILE:HD12	3:D:897:ILE:HA	1.80	0.42
1:A:3:ILE:HD12	1:A:3:ILE:HA	1.88	0.42
1:A:69:VAL:HG12	1:A:128:LEU:HG	2.01	0.42
3:D:363:PRO:HD3	5:F:52:TRP:CE2	2.54	0.42
3:D:1101:ASP:N	3:D:1101:ASP:OD1	2.51	0.42
7:H:17:DC:H2''	7:H:18:DA:C8	2.55	0.42
7:H:20:DG:H2''	7:H:21:DG:C8	2.55	0.41
1:B:124:HIS:HE1	1:B:127:THR:HG23	1.85	0.41
2:C:228:ARG:NH2	7:H:13:DC:H1'	2.35	0.41
2:C:563:ARG:HG2	2:C:564:LYS:H	1.84	0.41
3:D:876:ARG:HG2	3:D:1226:PHE:HZ	1.86	0.41
3:D:1089:PHE:HA	3:D:1095:SER:HA	2.03	0.41
2:C:58:THR:O	2:C:62:GLU:HG3	2.20	0.41
3:D:58:TRP:CE2	3:D:68:VAL:HG13	2.55	0.41
3:D:823:LEU:HD23	3:D:835:PRO:HB3	2.03	0.41
2:C:399:VAL:HG11	2:C:419:ASN:HB3	2.02	0.41
2:C:445:PRO:HD2	2:C:707:CYS:HB2	2.02	0.41
2:C:476:HIS:CG	2:C:477:PRO:HD2	2.55	0.41
3:D:1046:ILE:HD11	3:D:1120:GLU:HB3	2.01	0.41
7:H:21:DG:H2''	7:H:22:DA:C8	2.55	0.41
2:C:598:GLU:HB3	2:C:977:PHE:HD2	1.85	0.41
1:B:84:VAL:HG22	1:B:120:ASN:CG	2.46	0.41
1:B:202:ILE:HD13	1:B:207:ALA:HB2	2.02	0.41
2:C:454:ARG:HA	2:C:454:ARG:HD2	1.80	0.41
3:D:816:THR:HG23	3:D:821:LYS:HA	2.01	0.41
1:A:203:SER:OG	1:A:206:ASP:OD2	2.36	0.41
1:B:167:ILE:HG23	3:D:620:MET:HE1	2.03	0.41
2:C:454:ARG:NH1	2:C:499:SER:HB2	2.35	0.41
3:D:111:PRO:O	3:D:113:ARG:NH1	2.51	0.41
3:D:717:LYS:HE2	3:D:717:LYS:HB3	1.83	0.41
3:D:774:LEU:HD23	3:D:777:ILE:HD12	2.02	0.41
3:D:1152:LYS:HE3	3:D:1152:LYS:HB2	1.91	0.41
3:D:1152:LYS:O	3:D:1156:VAL:HG23	2.20	0.41
1:A:66:VAL:O	1:A:69:VAL:HG22	2.20	0.41
3:D:123:LYS:HB2	7:H:24:DG:OP1	2.21	0.41
1:B:42:LEU:HD21	1:B:208:LEU:HA	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:60:SER:HG	2:C:382:GLY:H	1.64	0.40
3:D:707:ILE:HD11	4:E:32:PRO:HB3	2.03	0.40
3:D:916:ILE:HG23	3:D:920:ALA:HB3	2.03	0.40
2:C:723:ILE:O	3:D:730:THR:HG23	2.21	0.40
4:E:33:LEU:HD23	4:E:33:LEU:H	1.85	0.40
5:F:131:ILE:H	5:F:131:ILE:HG13	1.69	0.40
3:D:1005:GLU:HB3	3:D:1006:PRO:HD3	2.03	0.40
1:A:83:LEU:HA	1:A:83:LEU:HD23	1.89	0.40
1:A:99:LYS:HG2	1:A:105:VAL:HG22	2.03	0.40
1:B:112:PRO:HA	1:B:113:PRO:HD2	1.96	0.40
3:D:789:LEU:HD22	3:D:793:TYR:CE2	2.53	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	223/359 (62%)	217 (97%)	5 (2%)	1 (0%)	30	64
1	B	230/359 (64%)	213 (93%)	16 (7%)	1 (0%)	30	64
2	C	1124/1178 (95%)	1072 (95%)	47 (4%)	5 (0%)	30	64
3	D	1261/1316 (96%)	1200 (95%)	58 (5%)	3 (0%)	43	74
4	E	79/110 (72%)	76 (96%)	3 (4%)	0	100	100
5	F	172/177 (97%)	168 (98%)	4 (2%)	0	100	100
All	All	3089/3499 (88%)	2946 (95%)	133 (4%)	10 (0%)	36	69

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	370	ILE

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Mol	Chain	Res	Type
1	B	159	ILE
2	C	732	GLU
1	A	184	GLU
2	C	32	VAL
3	D	593	PRO
3	D	1089	PHE
2	C	520	VAL
2	C	922	VAL
3	D	607	PRO

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	194/308 (63%)	187 (96%)	7 (4%)	31	54
1	B	191/308 (62%)	180 (94%)	11 (6%)	18	44
2	C	950/998 (95%)	925 (97%)	25 (3%)	40	60
3	D	1050/1095 (96%)	1030 (98%)	20 (2%)	50	66
4	E	66/90 (73%)	65 (98%)	1 (2%)	57	70
5	F	134/136 (98%)	130 (97%)	4 (3%)	36	57
All	All	2585/2935 (88%)	2517 (97%)	68 (3%)	40	60

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	18	ARG
1	A	51	VAL
1	A	59	VAL
1	A	111	VAL
1	A	116	VAL
1	A	171	VAL
1	B	17	ASN
1	B	79	ASN

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Mol	Chain	Res	Type
1	B	90	ASP
1	B	111	VAL
1	B	116	VAL
1	B	147	VAL
1	B	183	VAL
1	B	184	GLU
1	B	196	VAL
1	B	202	ILE
1	B	218	LEU
2	C	32	VAL
2	C	39	VAL
2	C	54	LEU
2	C	80	VAL
2	C	119	VAL
2	C	127	MET
2	C	185	VAL
2	C	208	ARG
2	C	224	VAL
2	C	228	ARG
2	C	246	GLU
2	C	322	LEU
2	C	363	VAL
2	C	370	ILE
2	C	373	PHE
2	C	409	VAL
2	C	571	VAL
2	C	593	MET
2	C	696	VAL
2	C	875	GLN
2	C	901	VAL
2	C	1057	LEU
2	C	1062	GLN
2	C	1137	VAL
2	C	1146	GLU
3	D	82	VAL
3	D	139	VAL
3	D	283	ASN
3	D	307	ASN
3	D	359	ASP
3	D	431	VAL
3	D	443	LEU
3	D	515	MET

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Mol	Chain	Res	Type
3	D	535	ASP
3	D	650	LEU
3	D	677	LEU
3	D	738	VAL
3	D	817	LEU
3	D	907	ASP
3	D	910	LEU
3	D	960	VAL
3	D	1009	GLN
3	D	1065	THR
3	D	1250	GLU
3	D	1279	ARG
4	E	84	GLU
5	F	56	GLU
5	F	58	ILE
5	F	60	ASP
5	F	86	PHE

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

Mol	Chain	Res	Type
2	C	142	ASN
2	C	375	ASN
2	C	386	GLN
2	C	419	ASN
2	C	543	GLN
2	C	614	GLN
2	C	685	ASN
2	C	700	GLN
2	C	755	HIS
2	C	918	ASN
2	C	1054	GLN
3	D	262	GLN
3	D	564	ASN
3	D	687	GLN
3	D	767	HIS
3	D	1109	GLN
5	F	101	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	225/359 (62%)	0.11	3 (1%) 75 54	16, 39, 97, 148	0
1	B	232/359 (64%)	0.25	8 (3%) 48 34	26, 63, 111, 150	0
2	C	1126/1178 (95%)	0.19	20 (1%) 67 47	12, 48, 141, 173	0
3	D	1265/1316 (96%)	0.21	17 (1%) 75 54	9, 61, 159, 205	0
4	E	81/110 (73%)	0.57	6 (7%) 20 19	49, 72, 145, 166	0
5	F	174/177 (98%)	0.60	12 (6%) 23 20	39, 86, 136, 162	0
6	G	10/10 (100%)	0.71	0 100 100	115, 135, 171, 178	0
7	H	24/24 (100%)	1.33	6 (25%) 2 4	125, 155, 191, 206	0
All	All	3137/3533 (88%)	0.24	72 (2%) 61 43	9, 58, 150, 206	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	F	78	ILE	5.3
2	C	237	LEU	4.5
2	C	32	VAL	3.8
2	C	252	PHE	3.8
3	D	1138	VAL	3.8
3	D	411	GLY	3.5
5	F	89	VAL	3.4
1	B	157	ALA	3.3
5	F	65	ALA	3.2
4	E	101	ALA	3.1
7	H	14	DT	3.1
3	D	827	PRO	3.1
2	C	248	ILE	3.0
3	D	854	HIS	3.0
5	F	4	VAL	2.9
3	D	416	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
2	C	249	VAL	2.8
1	B	156	GLY	2.8
2	C	1153	GLU	2.8
2	C	228	ARG	2.8
5	F	55	PRO	2.8
7	H	8	DA	2.7
4	E	108	GLU	2.7
2	C	568	VAL	2.7
3	D	412	ARG	2.7
4	E	28	GLY	2.6
3	D	825	THR	2.6
4	E	76	LEU	2.6
3	D	415	GLN	2.6
5	F	69	LEU	2.5
3	D	1098	VAL	2.5
7	H	10	DT	2.5
5	F	88	ASN	2.5
3	D	1280	ALA	2.5
7	H	15	DG	2.5
1	B	187	THR	2.4
2	C	828	LYS	2.4
3	D	1175	PHE	2.4
1	B	154	ALA	2.4
1	A	225	LEU	2.4
1	A	2	LEU	2.4
2	C	281	LEU	2.3
3	D	1142	TYR	2.3
2	C	28	SER	2.3
1	B	219	PHE	2.3
1	B	227	VAL	2.3
3	D	899	VAL	2.3
5	F	85	ARG	2.3
2	C	328	ILE	2.3
2	C	104	SER	2.2
5	F	80	GLU	2.2
2	C	560	LEU	2.2
7	H	25	DC	2.2
7	H	9	DG	2.2
1	B	62	GLU	2.2
2	C	584	ARG	2.2
1	B	2	LEU	2.2
2	C	566	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
4	E	68	TYR	2.1
1	A	3	ILE	2.1
5	F	99	PRO	2.1
2	C	229	LYS	2.1
3	D	8	ASP	2.1
3	D	956	GLY	2.1
3	D	1049	VAL	2.1
2	C	404	MET	2.1
5	F	177	ARG	2.1
3	D	1278	ALA	2.1
4	E	75	ILE	2.0
2	C	198	THR	2.0
2	C	240	ALA	2.0
5	F	84	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	ZN	D	2001	1/1	0.99	0.04	60,60,60,60	0
8	ZN	D	2002	1/1	1.00	0.02	68,68,68,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.