



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:18 AM UTC

PDB ID : 5E18 / pdb_00005e18
Title : T. thermophilus transcription initiation complex having a YYY discriminator sequence and a nontemplate-strand length corresponding to TSS selection at position 8 (RPo-CCC-8)
Authors : Zhang, Y.; Ebright, R.H.
Deposited on : 2015-09-29
Resolution : 3.30 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 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

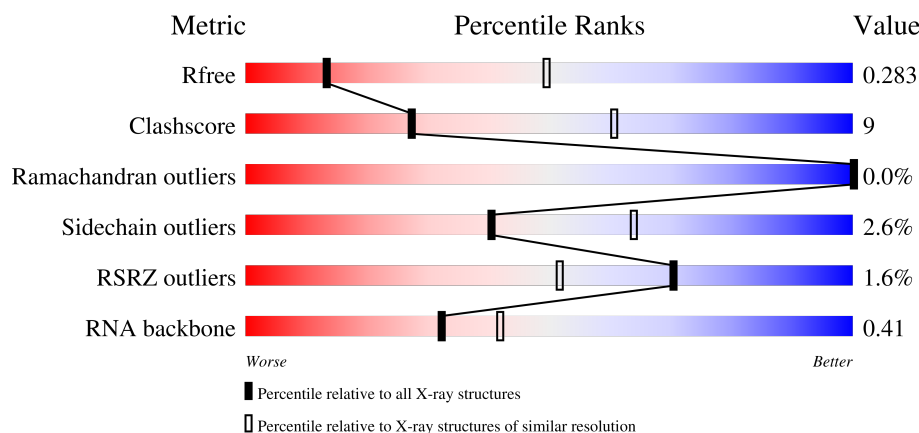
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)
RNA backbone	3983	1048 (3.60-3.00)

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	315	61% 12% 27%
1	B	315	52% 18% 29%
2	C	1119	75% 24% ..
3	D	1524	73% 24% ..

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Mol	Chain	Length	Quality of chain
4	E	99	2% 75% 18% • 5%
5	F	443	2% 59% 16% • 24%
6	G	21	38% 48% 14%
7	I	7	57% 29% 14%
8	H	28	43% 46% 11%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 28645 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	231	Total	C	N	O	S	0	1	0
			1814	1158	316	338	2			
1	B	223	Total	C	N	O	S	0	0	0
			1758	1124	305	327	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1112	Total	C	N	O	S	0	3	0
			8792	5562	1570	1636	24			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1486	Total	C	N	O	S	0	2	0
			11751	7450	2070	2195	36			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	94	Total	C	N	O	S	0	0	0
			758	483	132	139	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	335	Total	C	N	O	S	0	0	0
			2718	1713	497	504	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-19	MET	-	initiating methionine	UNP Q5SKW1
F	-18	GLY	-	expression tag	UNP Q5SKW1
F	-17	SER	-	expression tag	UNP Q5SKW1
F	-16	SER	-	expression tag	UNP Q5SKW1
F	-15	HIS	-	expression tag	UNP Q5SKW1
F	-14	HIS	-	expression tag	UNP Q5SKW1
F	-13	HIS	-	expression tag	UNP Q5SKW1
F	-12	HIS	-	expression tag	UNP Q5SKW1
F	-11	HIS	-	expression tag	UNP Q5SKW1
F	-10	HIS	-	expression tag	UNP Q5SKW1
F	-9	SER	-	expression tag	UNP Q5SKW1
F	-8	SER	-	expression tag	UNP Q5SKW1
F	-7	GLY	-	expression tag	UNP Q5SKW1
F	-6	LEU	-	expression tag	UNP Q5SKW1
F	-5	VAL	-	expression tag	UNP Q5SKW1
F	-4	PRO	-	expression tag	UNP Q5SKW1
F	-3	ARG	-	expression tag	UNP Q5SKW1
F	-2	GLY	-	expression tag	UNP Q5SKW1
F	-1	SER	-	expression tag	UNP Q5SKW1
F	0	HIS	-	expression tag	UNP Q5SKW1

- Molecule 6 is a DNA chain called DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*T
P*GP*AP*GP*TP*CP*GP*AP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	18	Total	C	N	O	P	0	0	0
			372	176	73	106	17			

- Molecule 7 is a RNA chain called RNA (5'-R(*CP*CP*CP*UP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	I	7	Total	C	N	O	P	0	0	0
			142	65	24	47	6			

- Molecule 8 is a DNA chain called DNA (28-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	25	Total	C	N	O	P	0	0	0
			508	243	93	148	24			

- Molecule 9 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total Mg 1 1	0	0
9	D	2	Total Mg 2 2	0	0
9	F	1	Total Mg 1 1	0	0

- Molecule 10 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

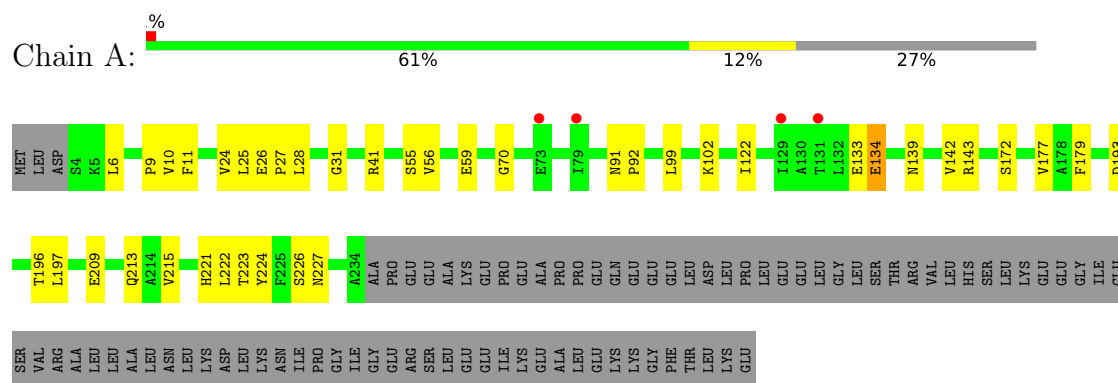
- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	1	Total O 1 1	0	0
11	B	3	Total O 3 3	0	0
11	C	8	Total O 8 8	0	0
11	D	10	Total O 10 10	0	0
11	E	1	Total O 1 1	0	0
11	F	2	Total O 2 2	0	0
11	H	1	Total O 1 1	0	0

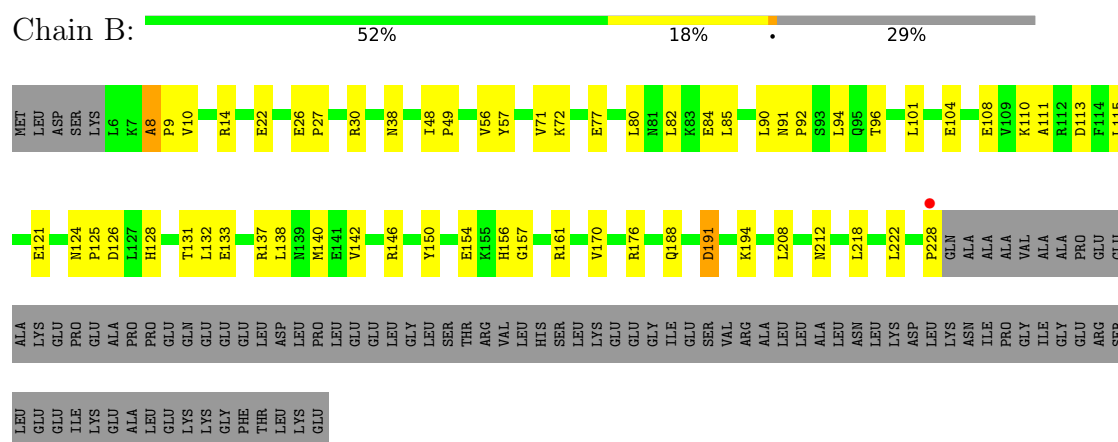
3 Residue-property plots

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

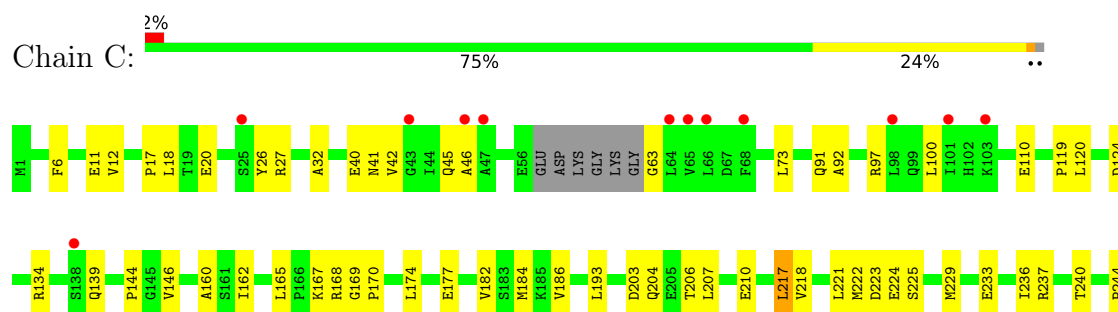
- Molecule 1: DNA-directed RNA polymerase subunit alpha



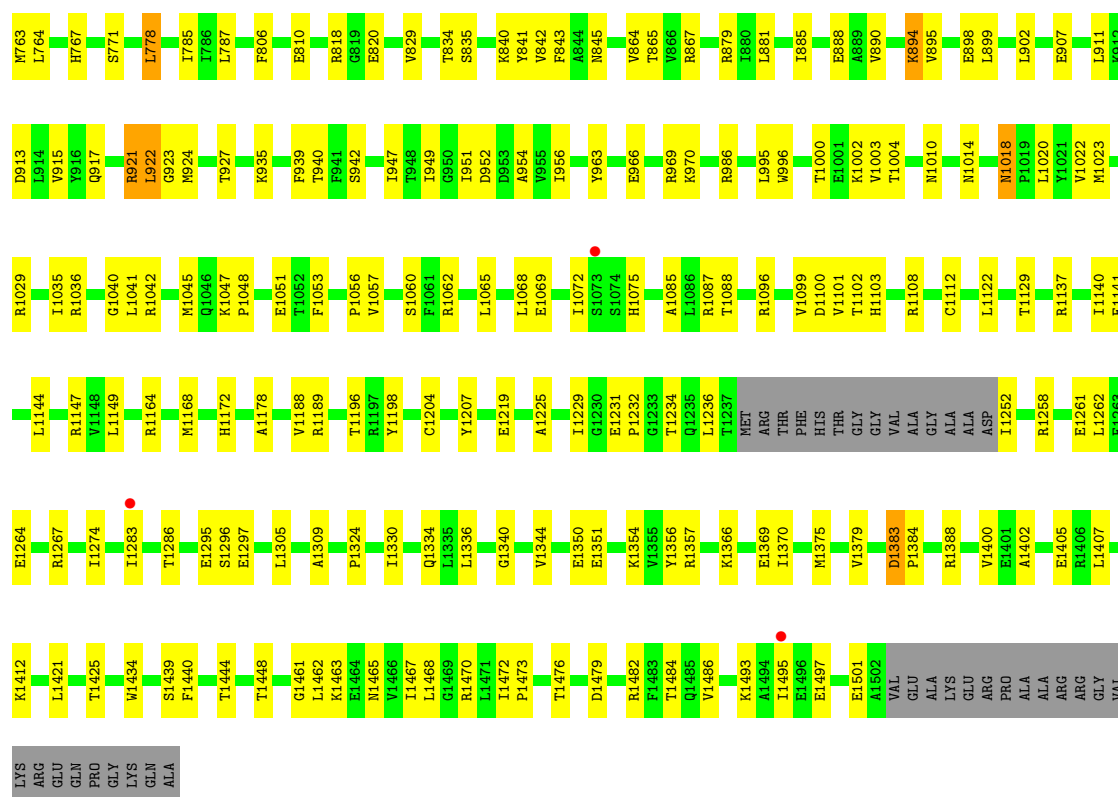
- Molecule 1: DNA-directed RNA polymerase subunit alpha



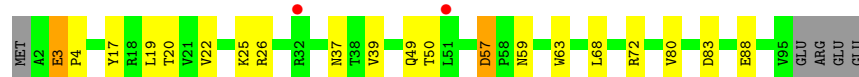
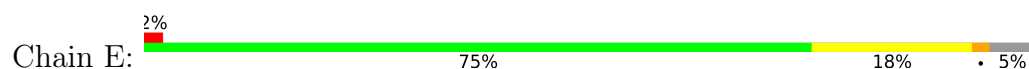
- Molecule 2: DNA-directed RNA polymerase subunit beta



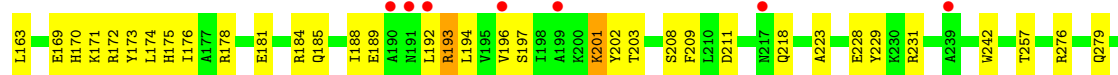
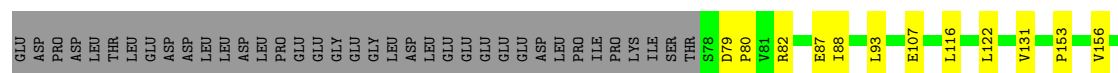
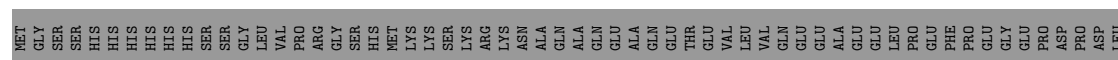




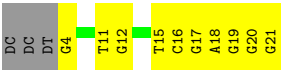
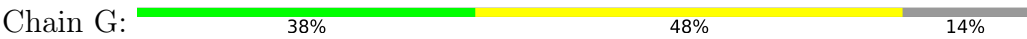
- Molecule 4: DNA-directed RNA polymerase subunit omega



- Molecule 5: RNA polymerase sigma factor SigA



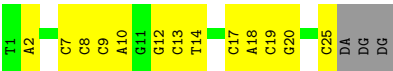
- Molecule 6: DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*CP*GP*AP*GP*GP*G)-3')



● Molecule 7: RNA (5'-R(*CP*CP*CP*UP*CP*GP*A)-3')



● Molecule 8: DNA (28-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.92Å 103.24Å 298.40Å 90.00° 97.93° 90.00°	Depositor
Resolution (Å)	38.34 – 3.30 38.34 – 3.30	Depositor EDS
% Data completeness (in resolution range)	90.3 (38.34-3.30) 91.3 (38.34-3.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.9	Depositor
R, R_{free}	0.238 , 0.284 0.242 , 0.283	Depositor DCC
R_{free} test set	3869 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	101.6	Xtriage
Anisotropy	0.584	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	28645	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	0.30	0/1849	0.81	4/2515 (0.2%)
1	B	0.31	0/1790	0.81	2/2435 (0.1%)
2	C	0.32	0/8969	0.77	7/12129 (0.1%)
3	D	0.34	0/11963	0.77	13/16173 (0.1%)
4	E	0.29	0/772	0.78	4/1040 (0.4%)
5	F	0.39	0/2759	0.81	3/3709 (0.1%)
6	G	0.20	0/418	0.30	0/645
7	I	0.14	0/157	0.29	0/242
8	H	0.29	0/569	0.48	0/876
All	All	0.33	0/29246	0.77	33/39764 (0.1%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	172	SER	CA-C-N	7.12	127.28	119.87
1	A	172	SER	C-N-CA	7.12	127.28	119.87
2	C	317	VAL	CA-C-N	6.97	127.01	119.90
2	C	317	VAL	C-N-CA	6.97	127.01	119.90
5	F	202	TYR	N-CA-C	6.20	120.45	113.02
3	D	1340	GLY	CA-C-N	6.07	125.28	118.97
3	D	1340	GLY	C-N-CA	6.07	125.28	118.97
2	C	876	VAL	CB-CA-C	-5.97	108.02	114.35
3	D	1383	ASP	CA-C-N	5.97	125.52	119.19
3	D	1383	ASP	C-N-CA	5.97	125.52	119.19
3	D	433	GLY	N-CA-C	5.87	119.32	110.75
1	B	115	LEU	CA-C-N	5.77	125.78	119.89
1	B	115	LEU	C-N-CA	5.77	125.78	119.89
2	C	980	GLY	N-CA-C	-5.71	106.64	114.64
5	F	333	ILE	CA-C-N	5.58	125.57	120.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	333	ILE	C-N-CA	5.58	125.57	120.21
3	D	1267	ARG	CA-C-N	5.45	125.46	119.90
3	D	1267	ARG	C-N-CA	5.45	125.46	119.90
2	C	439	CYS	CA-C-N	5.42	125.03	119.56
2	C	439	CYS	C-N-CA	5.42	125.03	119.56
4	E	57	ASP	CA-C-N	5.27	124.93	119.56
4	E	57	ASP	C-N-CA	5.27	124.93	119.56
3	D	1018	ASN	CA-C-N	5.21	125.26	119.32
3	D	1018	ASN	C-N-CA	5.21	125.26	119.32
1	A	227	ASN	CA-C-N	5.20	125.13	119.78
1	A	227	ASN	C-N-CA	5.20	125.13	119.78
3	D	845	ASN	CA-C-N	5.14	124.59	119.24
3	D	845	ASN	C-N-CA	5.14	124.59	119.24
2	C	303	PHE	N-CA-C	5.07	117.19	111.11
3	D	739	ASP	CA-C-N	-5.03	115.53	123.13
3	D	739	ASP	C-N-CA	-5.03	115.53	123.13
4	E	3	GLU	CA-C-N	5.00	125.28	119.93
4	E	3	GLU	C-N-CA	5.00	125.28	119.93

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	1814	0	1869	23	0
1	B	1758	0	1808	41	0
2	C	8792	0	8902	172	0
3	D	11751	0	11994	234	0
4	E	758	0	770	15	0
5	F	2718	0	2803	54	0
6	G	372	0	203	12	0
7	I	142	0	78	4	0
8	H	508	0	283	35	0
9	B	1	0	0	0	0
9	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	F	1	0	0	0	0
10	D	2	0	0	0	0
11	A	1	0	0	0	0
11	B	3	0	0	0	0
11	C	8	0	0	0	0
11	D	10	0	0	3	0
11	E	1	0	0	0	0
11	F	2	0	0	0	0
11	H	1	0	0	0	0
All	All	28645	0	28710	516	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 9.

All (516) 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)
6:G:20:DG:H1	7:I:2:C:H42	1.18	0.88
8:H:13:DC:O3'	8:H:14:DT:H72	1.74	0.86
5:F:209:PHE:HB2	8:H:9:DC:C4	2.12	0.84
2:C:628:PHE:H	2:C:638:ASP:HB3	1.43	0.83
2:C:950:LEU:HB3	2:C:952:LEU:HD13	1.66	0.77
2:C:167:LYS:HA	8:H:13:DC:C6	2.21	0.75
5:F:209:PHE:H	8:H:9:DC:N4	1.83	0.75
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.20	0.75
2:C:17:PRO:HB2	2:C:20:GLU:HB3	1.69	0.74
3:D:956:ILE:HD11	3:D:1062:ARG:HG2	1.71	0.72
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.23	0.71
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.73	0.70
3:D:132:TYR:OH	3:D:568:ARG:NH1	2.25	0.69
2:C:324:ASP:HB3	2:C:327:HIS:HB2	1.74	0.69
2:C:637:LEU:HG	2:C:659:PRO:HG3	1.73	0.69
3:D:65:ARG:NH1	5:F:378:GLY:O	2.27	0.67
8:H:13:DC:C2'	8:H:14:DT:H72	2.23	0.67
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.77	0.67
3:D:622:ARG:NH1	6:G:17:DG:OP1	2.27	0.67
3:D:657:LEU:HG	3:D:661:MET:HE2	1.77	0.66
3:D:1484:THR:O	4:E:25:LYS:NZ	2.21	0.66
5:F:209:PHE:H	8:H:9:DC:H42	1.41	0.66
2:C:1019:GLN:HG3	2:C:1058:ASP:HB3	1.75	0.66
1:B:14:ARG:HB3	1:B:22:GLU:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:939:ARG:HG2	2:C:982:PRO:HD3	1.77	0.65
2:C:244:PRO:O	5:F:82:ARG:NH1	2.29	0.65
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.78	0.65
3:D:63:TYR:HB2	3:D:80:VAL:HG21	1.79	0.65
2:C:168:ARG:HD3	2:C:268:ASP:HB3	1.78	0.65
3:D:241:ILE:HA	3:D:312:ARG:HG2	1.78	0.65
2:C:26:TYR:OH	2:C:119:PRO:O	2.13	0.64
2:C:711:GLU:O	2:C:758:ARG:NH1	2.29	0.64
5:F:194:LEU:O	5:F:197:SER:OG	2.14	0.64
8:H:9:DC:H2''	8:H:10:DA:C8	2.32	0.64
8:H:13:DC:O3'	8:H:14:DT:C7	2.44	0.64
2:C:768:THR:OG1	2:C:771:GLU:OE1	2.15	0.64
1:A:31:GLY:N	1:A:193:ASP:OD2	2.31	0.63
6:G:17:DG:H2'	6:G:18:DA:C8	2.34	0.62
2:C:146:VAL:HG22	2:C:162:ILE:HG12	1.81	0.62
1:B:77:GLU:OE1	3:D:867:ARG:NH2	2.32	0.62
3:D:140:ALA:HB2	3:D:452:ILE:HG12	1.81	0.61
1:B:26:GLU:HB3	1:B:194:LYS:HG3	1.82	0.61
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.82	0.61
3:D:1219:GLU:OE1	4:E:17:TYR:OH	2.17	0.61
2:C:12:VAL:HG21	2:C:472:ARG:HD3	1.83	0.61
2:C:428:ARG:NH2	2:C:447:ALA:O	2.32	0.61
3:D:566:ILE:HD11	5:F:192:LEU:HD21	1.81	0.61
6:G:21:DG:O6	7:I:1:C:N4	2.34	0.61
2:C:711:GLU:HG2	2:C:822:VAL:HG22	1.82	0.61
2:C:710:ILE:HD12	2:C:790:LEU:HB2	1.81	0.60
2:C:237:ARG:O	2:C:240:THR:OG1	2.18	0.60
2:C:343:GLN:NE2	2:C:384:GLU:OE2	2.33	0.60
3:D:192:ALA:HB3	3:D:195:VAL:HB	1.83	0.60
2:C:1100:GLN:HG3	3:D:9:ARG:HH21	1.66	0.60
3:D:685:ASP:HA	3:D:688:TRP:HD1	1.66	0.60
3:D:1096:ARG:NH1	3:D:1440:PHE:O	2.35	0.60
1:A:133:GLU:OE1	2:C:610:ARG:NH1	2.35	0.60
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.83	0.60
3:D:563:PRO:HD2	3:D:566:ILE:HD12	1.84	0.60
2:C:521:PRO:HB3	3:D:1068:LEU:HD21	1.84	0.59
2:C:805:ARG:HG2	2:C:807[A]:ARG:HE	1.66	0.59
2:C:727:PRO:HB3	2:C:783:ARG:HD3	1.84	0.59
3:D:1495:ILE:HD13	4:E:80:VAL:HG21	1.84	0.59
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.83	0.59
3:D:1479:ASP:OD1	3:D:1482:ARG:NE	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1486:VAL:HG11	4:E:26:ARG:HB2	1.84	0.58
2:C:167:LYS:HA	8:H:13:DC:H6	1.67	0.58
2:C:326:ASP:HA	2:C:331:ARG:HD2	1.86	0.58
3:D:643:GLY:HA3	3:D:727:GLN:HB2	1.84	0.58
1:B:104:GLU:OE2	1:B:137:ARG:NH1	2.37	0.58
3:D:256:GLU:OE2	3:D:300:LYS:NZ	2.27	0.58
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.85	0.58
3:D:970:LYS:HD3	3:D:995:LEU:HD13	1.84	0.58
3:D:218:LYS:HG2	3:D:338:GLU:HG2	1.86	0.58
3:D:894:LYS:H	3:D:894:LYS:HD3	1.68	0.58
1:B:38:ASN:ND2	2:C:978:ARG:O	2.35	0.57
3:D:1149:LEU:HD12	3:D:1164:ARG:HB3	1.87	0.57
2:C:946:ARG:NH1	11:D:2101:HOH:O	2.36	0.57
3:D:252:ARG:HA	3:D:303:PRO:HA	1.87	0.57
6:G:4:DG:H1	8:H:25:DC:H42	1.53	0.57
8:H:13:DC:C3'	8:H:14:DT:H72	2.35	0.57
1:B:191:ASP:OD1	1:B:191:ASP:N	2.33	0.57
3:D:715:ALA:HB3	3:D:764:LEU:HA	1.87	0.56
5:F:116:LEU:HD11	5:F:174:LEU:HA	1.86	0.56
1:A:215:VAL:HG13	1:B:222:LEU:HD22	1.87	0.56
2:C:223:ASP:OD2	2:C:225:SER:OG	2.21	0.56
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.87	0.56
2:C:674:VAL:HG12	2:C:869:VAL:HB	1.88	0.56
2:C:97:ARG:NH1	2:C:110:GLU:OE1	2.36	0.56
3:D:129:PHE:CD2	3:D:456:MET:HB3	2.41	0.56
2:C:399:ASN:O	2:C:402:SER:OG	2.20	0.56
2:C:502:PRO:HB2	2:C:509:ALA:HB3	1.88	0.56
3:D:162:ARG:O	3:D:414:ARG:NH1	2.37	0.56
8:H:13:DC:H2''	8:H:14:DT:H72	1.87	0.56
2:C:884:GLN:O	2:C:888:THR:OG1	2.23	0.56
3:D:703:ASN:HB2	3:D:713:ILE:HG12	1.88	0.56
5:F:209:PHE:CB	8:H:9:DC:C4	2.87	0.56
2:C:545:ASN:HB3	2:C:583:LEU:HD22	1.87	0.55
2:C:168:ARG:O	2:C:267:TYR:HA	2.06	0.55
5:F:209:PHE:N	8:H:9:DC:N4	2.55	0.55
2:C:971:LYS:HG2	2:C:988:VAL:HG22	1.88	0.55
2:C:167:LYS:HD3	8:H:13:DC:H2'	1.88	0.55
2:C:1019:GLN:OE1	3:D:617:ASN:ND2	2.29	0.55
2:C:456:ALA:HB3	2:C:459:ALA:HB2	1.88	0.55
3:D:433:GLY:HA2	3:D:449:SER:H	1.72	0.55
3:D:1147:ARG:NH2	3:D:1369:GLU:OE1	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:627:ARG:HD3	2:C:638:ASP:HB2	1.89	0.55
3:D:1003:VAL:HG21	3:D:1041:LEU:HG	1.89	0.54
1:B:84:GLU:OE2	3:D:867:ARG:NH1	2.37	0.54
3:D:1468:LEU:HD23	3:D:1470:ARG:HD2	1.88	0.54
2:C:1016:ILE:O	3:D:87:ARG:NH1	2.38	0.54
2:C:1036:GLU:OE2	2:C:1036:GLU:N	2.40	0.54
1:A:223:THR:O	1:A:226:SER:OG	2.21	0.54
1:B:30:ARG:HH21	2:C:854:PRO:HB3	1.73	0.54
5:F:284:ARG:NH2	5:F:290:GLU:OE2	2.41	0.54
2:C:503:LEU:HD23	2:C:508:ILE:HA	1.89	0.54
3:D:1375:MET:HE2	3:D:1421:LEU:HD11	1.89	0.54
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.90	0.54
3:D:272:LEU:HB2	3:D:280:ALA:HB3	1.90	0.54
3:D:741:ASP:OD1	3:D:741:ASP:N	2.40	0.54
5:F:171:LYS:O	5:F:175:HIS:ND1	2.30	0.54
1:B:138:LEU:HD11	1:B:140:MET:HE3	1.89	0.54
2:C:167:LYS:CA	8:H:13:DC:C6	2.90	0.54
2:C:915:LYS:NZ	3:D:952:ASP:OD2	2.40	0.54
5:F:209:PHE:N	8:H:9:DC:H42	2.05	0.54
2:C:614:ARG:NH2	2:C:618:GLY:O	2.41	0.53
3:D:785:ILE:HD13	3:D:935:LYS:HA	1.89	0.53
3:D:181:ASP:HB2	3:D:205:TYR:CD2	2.43	0.53
2:C:937:ASP:OD1	2:C:939:ARG:HD3	2.08	0.53
3:D:1462:LEU:HD23	3:D:1473:PRO:HD2	1.89	0.53
2:C:405:ARG:HD3	2:C:566:THR:HG21	1.89	0.53
2:C:1083:GLU:OE2	3:D:87:ARG:NH2	2.40	0.53
3:D:1264:GLU:OE2	3:D:1425:THR:OG1	2.26	0.53
1:A:133:GLU:HG2	1:A:134:GLU:H	1.73	0.53
1:A:209:GLU:O	1:A:213:GLN:HG2	2.08	0.53
2:C:396:ASP:HA	2:C:633:GLN:NE2	2.23	0.53
2:C:550:LEU:HD23	2:C:906:PHE:HE1	1.74	0.53
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.42	0.53
3:D:114:THR:HG23	3:D:495:ARG:HG2	1.90	0.52
3:D:435:VAL:HG22	3:D:446:VAL:HG22	1.91	0.52
3:D:1085:ALA:O	3:D:1088:THR:OG1	2.22	0.52
1:B:111:ALA:HB3	1:B:125:PRO:HA	1.89	0.52
6:G:15:DT:H2'	6:G:16:DC:C6	2.44	0.52
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.91	0.52
8:H:12:DG:C8	8:H:12:DG:H5''	2.45	0.52
2:C:92:ALA:HB2	2:C:120:LEU:HD11	1.91	0.52
2:C:657:ASP:OD2	2:C:663:ASN:N	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:127:LEU:HA	3:D:457:GLY:HA2	1.92	0.52
3:D:208:PRO:HA	3:D:390:PRO:HA	1.91	0.52
3:D:840:LYS:HE3	3:D:841:TYR:CZ	2.45	0.52
2:C:203:ASP:OD1	2:C:204:GLN:N	2.43	0.52
3:D:12:LEU:HD21	3:D:104:PHE:CZ	2.44	0.52
3:D:834:THR:OG1	3:D:835:SER:N	2.42	0.52
5:F:365:GLU:HB2	5:F:404:ALA:HB2	1.92	0.52
3:D:761:ILE:HD12	4:E:20:THR:HA	1.92	0.52
3:D:1047:LYS:HD3	3:D:1051:GLU:HB3	1.91	0.52
3:D:1350:GLU:O	3:D:1354:LYS:HG3	2.10	0.52
1:A:59:GLU:OE1	1:A:139[B]:ASN:ND2	2.30	0.51
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.91	0.51
5:F:193:ARG:HG2	8:H:7:DC:H5''	1.93	0.51
2:C:617:ASP:OD1	2:C:617:ASP:N	2.43	0.51
2:C:705:ILE:HA	2:C:828:ALA:HA	1.91	0.51
2:C:495:THR:N	2:C:530:GLU:OE1	2.42	0.51
2:C:1092:LEU:HD13	2:C:1099:VAL:HG21	1.93	0.51
3:D:97:THR:HG21	3:D:571:LYS:HG2	1.92	0.51
3:D:996:TRP:CD2	3:D:1056:PRO:HG3	2.45	0.51
2:C:669:GLY:HA3	2:C:995:MET:HA	1.93	0.51
3:D:25:GLU:HB2	3:D:92:HIS:CE1	2.46	0.51
3:D:187:LYS:N	3:D:200:ASP:OD2	2.34	0.51
3:D:273:ARG:HB3	3:D:278:PRO:HA	1.93	0.51
3:D:314:PRO:HB2	3:D:317:VAL:HG12	1.93	0.51
3:D:1324:PRO:HG3	3:D:1330:ILE:HD11	1.93	0.51
3:D:1305:LEU:HD13	3:D:1309:ALA:HB3	1.92	0.50
6:G:18:DA:H2'	6:G:19:DG:C8	2.45	0.50
1:A:70:GLY:H	2:C:607:ASP:CG	2.19	0.50
2:C:1038:TRP:CE2	3:D:1099:VAL:HG11	2.47	0.50
3:D:1004:THR:HG23	3:D:1036:ARG:HB2	1.94	0.50
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.12	0.50
8:H:13:DC:H1'	8:H:14:DT:C7	2.41	0.50
1:A:55:SER:HB3	1:A:143:ARG:HB3	1.93	0.50
1:B:8:ALA:HB1	1:B:9:PRO:HA	1.93	0.50
2:C:374:ASN:OD1	5:F:276:ARG:HD2	2.12	0.50
3:D:58:CYS:SG	3:D:62:LYS:N	2.85	0.50
3:D:890:VAL:HB	3:D:922:LEU:HD12	1.93	0.50
1:A:11:PHE:O	1:B:228:PRO:HA	2.11	0.49
3:D:132:TYR:HB2	3:D:153:LEU:HB2	1.94	0.49
4:E:37:ASN:OD1	4:E:37:ASN:N	2.42	0.49
5:F:208:SER:OG	8:H:9:DC:N4	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:551:GLU:OE2	2:C:551:GLU:N	2.35	0.49
3:D:534:ARG:HH12	5:F:312:GLN:HB3	1.77	0.49
5:F:189:GLU:O	5:F:192:LEU:HB2	2.11	0.49
5:F:228:GLU:OE1	5:F:231:ARG:NH2	2.31	0.49
2:C:1110:ASP:OD2	2:C:1114:GLY:N	2.37	0.49
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.12	0.49
3:D:1232:PRO:HG2	3:D:1356:TYR:HE1	1.77	0.49
2:C:224:GLU:CD	2:C:224:GLU:H	2.20	0.49
2:C:704:HIS:CD2	2:C:831:ARG:HD2	2.47	0.49
3:D:1065:LEU:HD23	3:D:1069:GLU:HB3	1.93	0.49
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.93	0.49
3:D:256:GLU:OE1	3:D:274:ARG:NH1	2.45	0.49
3:D:472:ALA:O	3:D:476:GLU:HG2	2.12	0.49
3:D:1042:ARG:HB3	3:D:1057:VAL:HB	1.93	0.49
3:D:100:ALA:HB3	3:D:575:GLN:HE22	1.78	0.49
3:D:633:VAL:HB	3:D:740:PHE:CZ	2.48	0.49
3:D:1020:LEU:HB3	3:D:1035:ILE:HD12	1.94	0.49
3:D:1366:LYS:O	3:D:1370:ILE:HG12	2.12	0.49
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.94	0.49
3:D:1101:VAL:HG13	3:D:1102:THR:HG23	1.95	0.49
5:F:203:THR:HB	8:H:9:DC:O2	2.13	0.49
2:C:361:MET:HE2	5:F:201:LYS:HD2	1.94	0.49
2:C:471:TYR:OH	2:C:516:ARG:NH2	2.45	0.49
1:B:56:VAL:HG21	1:B:82:LEU:HD13	1.95	0.48
1:B:30:ARG:NH2	2:C:854:PRO:HB3	2.28	0.48
2:C:496:ILE:HG12	2:C:531:PHE:HB2	1.95	0.48
2:C:751:PRO:HB3	2:C:794:PRO:HA	1.95	0.48
3:D:898:GLU:OE2	3:D:921:ARG:NH2	2.46	0.48
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.94	0.48
3:D:1168:MET:HE2	3:D:1172:HIS:CE1	2.48	0.48
2:C:704:HIS:CE1	2:C:1000:MET:HE1	2.48	0.48
5:F:276:ARG:O	5:F:279:GLN:HG3	2.12	0.48
1:B:150:TYR:CE2	1:B:170:VAL:HG22	2.49	0.48
5:F:170:HIS:HA	5:F:173:TYR:HD2	1.78	0.48
5:F:181:GLU:O	5:F:184:ARG:HB3	2.13	0.48
8:H:8:DC:H2"	8:H:9:DC:OP2	2.14	0.48
1:B:176:ARG:NH2	3:D:888:GLU:OE1	2.46	0.48
2:C:170:PRO:HA	8:H:14:DT:O4	2.14	0.48
2:C:304:LEU:HB3	2:C:305:PRO:HD3	1.95	0.48
5:F:172:ARG:O	5:F:176:ILE:HG12	2.14	0.48
3:D:778:LEU:HD12	3:D:778:LEU:HA	1.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:613:ARG:HG3	3:D:618:LEU:HD23	1.96	0.48
2:C:124:ASP:OD2	2:C:407:LYS:NZ	2.47	0.48
5:F:223:ALA:HB2	5:F:242:TRP:HB2	1.97	0.47
2:C:715:THR:OG1	2:C:718:GLY:O	2.28	0.47
3:D:96:ALA:HB2	3:D:555:LYS:HG2	1.95	0.47
4:E:57:ASP:O	4:E:63:TRP:NE1	2.41	0.47
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.75	0.47
2:C:571:LEU:HB2	2:C:574:ALA:HB2	1.96	0.47
1:A:25:LEU:HD23	1:A:28:LEU:HD11	1.96	0.47
2:C:42:VAL:O	2:C:45:GLN:HB3	2.13	0.47
2:C:396:ASP:HA	2:C:633:GLN:HE22	1.80	0.47
3:D:1189:ARG:HB3	3:D:1204:CYS:HA	1.96	0.47
4:E:39:VAL:O	4:E:72:ARG:NH1	2.42	0.47
8:H:19:DC:H2"	8:H:20:DG:C8	2.49	0.47
5:F:395:GLU:OE2	5:F:398:ARG:NH2	2.48	0.47
2:C:537:LYS:HD3	2:C:583:LEU:HD11	1.96	0.47
3:D:171:LEU:HD12	3:D:390:PRO:HG2	1.96	0.47
2:C:549:PHE:HB3	2:C:552:HIS:HD2	1.78	0.47
2:C:881:ASN:OD1	2:C:881:ASN:N	2.48	0.47
3:D:42:ASP:N	3:D:46:ASP:OD2	2.38	0.47
3:D:1379:VAL:HG21	3:D:1400:VAL:HG11	1.97	0.47
1:B:208:LEU:O	1:B:212:ASN:ND2	2.48	0.47
2:C:704:HIS:O	2:C:829:GLN:HG2	2.15	0.47
2:C:18:LEU:HD13	2:C:404:LEU:HD21	1.97	0.47
3:D:142:LEU:HB2	3:D:161:LEU:HD11	1.97	0.47
1:B:80:LEU:HD21	3:D:842:VAL:HG12	1.96	0.46
2:C:41:ASN:O	2:C:46:ALA:HB2	2.16	0.46
3:D:455:ARG:HB2	3:D:460:ALA:HB2	1.97	0.46
5:F:169:GLU:O	5:F:172:ARG:HB3	2.16	0.46
1:A:221:HIS:HA	1:A:224:TYR:CD2	2.50	0.46
2:C:437:ARG:NH2	2:C:491:GLU:OE2	2.47	0.46
3:D:619:LEU:HD11	3:D:1439:SER:HB2	1.97	0.46
3:D:1112:CYS:HB3	3:D:1196:THR:OG1	2.15	0.46
2:C:930:LYS:HE3	2:C:935:GLY:HA2	1.98	0.46
2:C:1019:GLN:CG	2:C:1058:ASP:HB3	2.45	0.46
1:A:99:LEU:HD21	1:A:122:ILE:HD11	1.98	0.46
1:B:124:ASN:OD1	1:B:124:ASN:N	2.48	0.46
3:D:178:LEU:HG	3:D:192:ALA:HA	1.98	0.46
3:D:452:ILE:H	3:D:452:ILE:HG13	1.36	0.46
3:D:761:ILE:HD13	4:E:19:LEU:HG	1.97	0.46
3:D:1402:ALA:O	3:D:1405:GLU:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:PRO:HB3	1:A:27:PRO:O	2.15	0.46
2:C:740:GLU:HB3	2:C:805:ARG:HH12	1.81	0.46
3:D:1258:ARG:NH2	3:D:1262:LEU:HD21	2.30	0.46
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.97	0.46
2:C:1042:ALA:HB3	3:D:710:ARG:HB3	1.98	0.46
3:D:1045[B]:MET:HE3	3:D:1045[B]:MET:HB2	1.58	0.46
3:D:1486:VAL:HG22	4:E:22:VAL:HG13	1.98	0.46
1:A:102:LYS:HE3	1:A:102:LYS:HB2	1.74	0.46
3:D:500:ARG:NH1	3:D:1388:ARG:O	2.47	0.46
3:D:758:GLU:OE1	3:D:1476:THR:OG1	2.26	0.46
2:C:874:LEU:HB3	3:D:1029:ARG:HG3	1.98	0.46
2:C:893:ALA:HB2	2:C:918:LEU:HD23	1.98	0.46
3:D:658:LEU:HA	3:D:661:MET:HE3	1.98	0.46
3:D:1225:ALA:O	3:D:1229:ILE:HG13	2.16	0.46
3:D:106:LYS:O	3:D:586:ARG:NH1	2.50	0.45
3:D:483:HIS:CG	3:D:484:PRO:HD2	2.51	0.45
1:B:57:TYR:CG	1:B:161:ARG:HD2	2.51	0.45
2:C:1043:TYR:OH	3:D:713:ILE:O	2.28	0.45
3:D:411:THR:HG23	3:D:436:GLU:HA	1.98	0.45
6:G:20:DG:N2	7:I:2:C:N3	2.46	0.45
2:C:1005:MET:HE2	3:D:724:GLN:HE21	1.82	0.45
3:D:468:LEU:HD23	3:D:468:LEU:HA	1.83	0.45
3:D:633:VAL:O	3:D:635:PRO:HD3	2.16	0.45
5:F:208:SER:HB3	5:F:211:ASP:OD2	2.17	0.45
5:F:333:ILE:HA	5:F:334:PRO:HD3	1.82	0.45
1:B:14:ARG:O	1:B:22:GLU:N	2.45	0.45
2:C:612:VAL:HG22	2:C:622:GLU:HG3	1.97	0.45
2:C:1019:GLN:HE21	3:D:621:LYS:HE3	1.80	0.45
3:D:963:TYR:CE1	3:D:1002:LYS:HD3	2.52	0.45
3:D:1383:ASP:HA	3:D:1384:PRO:HD3	1.75	0.45
8:H:13:DC:O3'	8:H:14:DT:C5	2.69	0.45
2:C:1043:TYR:CD2	3:D:763:MET:HG2	2.52	0.45
3:D:923:GLY:O	3:D:927:THR:OG1	2.25	0.45
3:D:1258:ARG:NH1	3:D:1261:GLU:OE2	2.47	0.45
2:C:470:PRO:HB2	2:C:534:VAL:HG21	1.99	0.45
3:D:924:MET:HE2	3:D:924:MET:HB2	1.70	0.45
3:D:1296:SER:OG	3:D:1297:GLU:N	2.50	0.45
3:D:1493:LYS:O	3:D:1497:GLU:HG2	2.17	0.45
5:F:285:GLU:HA	5:F:286:PRO:HD3	1.82	0.45
2:C:139:GLN:HB2	2:C:391:LEU:HD11	1.99	0.45
2:C:553:ASP:OD2	2:C:843:HIS:ND1	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:286:SER:OG	2:C:301:GLU:OE2	2.26	0.45
2:C:317:VAL:O	2:C:320:HIS:ND1	2.42	0.45
2:C:261:ILE:HG22	2:C:262:ALA:N	2.32	0.44
3:D:939:PHE:O	3:D:942:SER:OG	2.29	0.44
2:C:134:ARG:NH2	6:G:21:DG:O3'	2.35	0.44
2:C:615:TYR:OH	2:C:623:TYR:OH	2.15	0.44
3:D:584:ASN:HD21	3:D:591:VAL:H	1.65	0.44
3:D:764:LEU:HD23	3:D:767:HIS:CD2	2.53	0.44
3:D:1122:LEU:HD13	3:D:1178:ALA:HB2	1.99	0.44
3:D:1137:ARG:O	3:D:1141:GLU:HG3	2.17	0.44
2:C:712:ALA:O	2:C:820:ARG:N	2.49	0.44
3:D:881:LEU:O	3:D:885:ILE:HG13	2.17	0.44
2:C:356:ARG:HA	2:C:359:MET:HE2	1.99	0.44
2:C:420:ARG:C	2:C:422:ARG:H	2.25	0.44
3:D:684:LYS:HB3	3:D:684:LYS:HE2	1.78	0.44
3:D:954:ALA:O	3:D:1062:ARG:NH2	2.50	0.44
3:D:1351:GLU:O	3:D:1354:LYS:HB2	2.16	0.44
3:D:1444:THR:O	3:D:1448:THR:HG23	2.17	0.44
2:C:578:VAL:HA	2:C:900:ARG:HG2	2.00	0.44
3:D:101:HIS:CE1	3:D:582:LEU:HD13	2.52	0.44
3:D:644:LEU:HD12	3:D:645:PRO:HD2	2.00	0.44
3:D:1047:LYS:HG2	3:D:1053:PHE:CE2	2.52	0.44
5:F:364:ARG:HH12	5:F:396:ARG:NH2	2.15	0.44
1:B:110:LYS:HD3	1:B:128:HIS:HA	2.00	0.44
2:C:343:GLN:HG3	2:C:385:PHE:CD1	2.53	0.44
2:C:390:GLN:HB3	2:C:415:PRO:HD3	2.00	0.44
2:C:1101:THR:OG1	2:C:1109:VAL:O	2.33	0.44
3:D:1100:ASP:OD2	3:D:1463:LYS:NZ	2.30	0.44
5:F:361:LEU:HD11	5:F:408:LEU:HG	1.99	0.44
8:H:12:DG:H4'	8:H:13:DC:OP2	2.17	0.44
1:A:56:VAL:HG22	1:A:142:VAL:HG12	1.99	0.44
1:B:128:HIS:CE1	1:B:131:THR:HG23	2.52	0.44
2:C:1031:ARG:HA	3:D:622:ARG:HA	1.99	0.44
3:D:65:ARG:HA	3:D:65:ARG:HD3	1.69	0.44
3:D:806:PHE:HB2	3:D:829:VAL:HG22	2.00	0.44
5:F:153:PRO:HA	5:F:156:VAL:HG22	2.00	0.44
3:D:966:GLU:O	3:D:969:ARG:HG2	2.18	0.44
3:D:1274:ILE:HD11	3:D:1334:GLN:HB3	2.00	0.44
1:B:156:HIS:CG	1:B:157:GLY:H	2.35	0.43
2:C:193:LEU:HD23	2:C:307:LEU:HD22	2.00	0.43
2:C:551:GLU:HB2	3:D:1065:LEU:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:487:ALA:O	3:D:491:LYS:HG2	2.18	0.43
3:D:913:ASP:O	3:D:917:GLN:HG2	2.18	0.43
3:D:1010:ASN:OD1	3:D:1014:ASN:ND2	2.46	0.43
2:C:1048:THR:O	2:C:1052:MET:HG2	2.18	0.43
3:D:31:THR:HG21	5:F:257:THR:HG22	2.00	0.43
3:D:96:ALA:HB3	3:D:554:LEU:HD23	2.00	0.43
3:D:1048:PRO:HD3	3:D:1075:HIS:CD2	2.54	0.43
2:C:160:ALA:HB3	2:C:174:LEU:HB2	2.00	0.43
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.51	0.43
3:D:230:TRP:CZ3	3:D:232:GLU:HG2	2.53	0.43
3:D:1462:LEU:HD22	3:D:1472:ILE:HB	1.99	0.43
5:F:107:GLU:HG3	5:F:229:TYR:HD2	1.83	0.43
6:G:11:DT:H2"	6:G:12:DG:H5"	2.00	0.43
3:D:648:MET:O	3:D:652:LEU:HG	2.18	0.43
3:D:911:LEU:O	3:D:915:VAL:HG23	2.19	0.43
6:G:4:DG:H1	8:H:25:DC:N4	2.13	0.43
2:C:91:GLN:HB3	2:C:119:PRO:HA	2.00	0.43
2:C:217:LEU:HD13	2:C:217:LEU:HA	1.82	0.43
2:C:458:TYR:HB3	2:C:470:PRO:HG3	2.01	0.43
2:C:922:PHE:CE2	2:C:964:LYS:HB2	2.53	0.43
3:D:63:TYR:HE2	3:D:73:CYS:HA	1.82	0.43
3:D:543:LEU:HD13	3:D:581:LEU:HA	2.01	0.43
3:D:879:ARG:HD3	3:D:902:LEU:O	2.18	0.43
1:B:72:LYS:HB3	1:B:131:THR:OG1	2.18	0.43
1:B:90:LEU:HD21	1:B:121:GLU:HB2	2.01	0.43
2:C:859:PRO:O	2:C:867:VAL:HG22	2.19	0.43
2:C:1056:LYS:HB3	3:D:623:VAL:HB	2.00	0.43
2:C:1097:LEU:HD11	3:D:103:TRP:HZ3	1.83	0.43
3:D:356:PRO:HB3	3:D:441:ARG:HA	2.00	0.43
3:D:949:ILE:HD11	3:D:1023:MET:HE1	2.00	0.43
1:B:8:ALA:HB1	1:B:27:PRO:HD2	2.00	0.43
2:C:40:GLU:O	2:C:45:GLN:HG2	2.18	0.43
3:D:125:GLN:HA	3:D:130:SER:HB3	2.01	0.43
3:D:1140:ILE:O	3:D:1144:LEU:HB2	2.18	0.43
2:C:139:GLN:NE2	2:C:413:LEU:O	2.52	0.43
2:C:891:GLY:O	2:C:991:GLN:HB2	2.18	0.43
3:D:523:ASP:O	3:D:526:PRO:HG3	2.19	0.43
3:D:671:LYS:NZ	5:F:421:PHE:HA	2.34	0.43
1:B:156:HIS:CG	1:B:157:GLY:N	2.86	0.43
2:C:578:VAL:HG23	2:C:671:ASN:CG	2.44	0.43
2:C:1054:THR:OG1	2:C:1055:LEU:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:618:LEU:HD12	3:D:1467:ILE:HG23	1.99	0.43
5:F:196:VAL:HG11	8:H:7:DC:H4'	2.01	0.43
3:D:771:SER:HB2	3:D:778:LEU:HD22	2.00	0.43
3:D:843:PHE:HE2	3:D:864:VAL:HG21	1.84	0.43
1:B:128:HIS:HE1	1:B:131:THR:HG23	1.83	0.42
2:C:32:ALA:HB2	2:C:73:LEU:HD12	2.01	0.42
1:A:221:HIS:O	1:A:224:TYR:HB2	2.20	0.42
2:C:419:THR:HG23	2:C:422:ARG:HG3	2.01	0.42
3:D:161:LEU:HB3	3:D:452:ILE:HD11	2.01	0.42
6:G:17:DG:H2'	6:G:18:DA:H8	1.80	0.42
8:H:12:DG:H3'	8:H:13:DC:O4'	2.19	0.42
2:C:218:VAL:HG22	2:C:222:MET:HE2	2.01	0.42
2:C:543:ASN:HD21	2:C:566:THR:HG22	1.84	0.42
2:C:627:ARG:HH22	2:C:640:ARG:HG3	1.84	0.42
3:D:787:LEU:HD21	3:D:947:ILE:HG21	2.02	0.42
3:D:1463:LYS:HE2	3:D:1463:LYS:HB3	1.78	0.42
2:C:236:ILE:O	2:C:240:THR:HG23	2.20	0.42
2:C:595:LEU:HB3	2:C:655:LEU:HB2	2.02	0.42
2:C:1038:TRP:NE1	3:D:1099:VAL:HG11	2.35	0.42
3:D:8:VAL:HG12	3:D:1434:TRP:HZ2	1.85	0.42
3:D:1144:LEU:HD23	3:D:1144:LEU:HA	1.76	0.42
5:F:107:GLU:HG3	5:F:229:TYR:CD2	2.54	0.42
2:C:144:PRO:HG2	2:C:165:LEU:HD23	2.01	0.42
2:C:299:LYS:HB2	2:C:299:LYS:HE3	1.84	0.42
3:D:701:LEU:HD13	3:D:763:MET:HE1	2.02	0.42
3:D:1072:ILE:O	3:D:1075:HIS:HB2	2.20	0.42
3:D:1087:ARG:HD2	3:D:1236:LEU:O	2.19	0.42
5:F:163:LEU:HB3	5:F:174:LEU:HD22	2.02	0.42
5:F:318:GLU:OE1	5:F:318:GLU:N	2.50	0.42
2:C:888:THR:O	2:C:990:GLY:HA3	2.20	0.42
2:C:910:LYS:O	2:C:914:ILE:HG13	2.19	0.42
3:D:200:ASP:O	3:D:397:LYS:HG2	2.20	0.42
3:D:572:ARG:NH2	5:F:87:GLU:OE2	2.53	0.42
3:D:818:ARG:HE	3:D:820:GLU:CD	2.27	0.42
1:B:48:ILE:HA	1:B:49:PRO:HD3	1.89	0.42
1:B:85:LEU:HA	1:B:124:ASN:HD21	1.85	0.42
2:C:389:SER:OG	2:C:390:GLN:N	2.53	0.42
2:C:390:GLN:HE21	2:C:390:GLN:HB2	1.57	0.42
3:D:15:PRO:HG3	11:D:2105:HOH:O	2.20	0.42
3:D:258:VAL:HG12	3:D:273:ARG:O	2.20	0.42
3:D:840:LYS:HE3	3:D:841:TYR:OH	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:184:ARG:O	5:F:188:ILE:HG13	2.20	0.42
2:C:944:LEU:HD21	2:C:963:LEU:HD23	2.02	0.42
2:C:992:MET:HG2	2:C:994:ILE:HG13	2.02	0.42
3:D:689:ASP:O	3:D:693:GLU:HG3	2.19	0.42
3:D:1147:ARG:HH22	3:D:1369:GLU:CD	2.27	0.42
2:C:571:LEU:HD22	2:C:700:TYR:HA	2.02	0.42
2:C:678:PRO:HA	2:C:683:ASN:HD21	1.84	0.42
5:F:181:GLU:O	5:F:185:GLN:HG2	2.20	0.42
1:B:91:ASN:HA	1:B:92:PRO:HD2	1.84	0.41
4:E:3:GLU:HA	4:E:4:PRO:HD3	1.84	0.41
5:F:231:ARG:HG3	8:H:2:DA:C2	2.54	0.41
8:H:13:DC:H1'	8:H:14:DT:H72	2.02	0.41
1:A:10:VAL:HG12	1:A:26:GLU:O	2.19	0.41
1:A:41:ARG:HA	1:A:177:VAL:HG11	2.01	0.41
3:D:648:MET:HE2	3:D:648:MET:HB3	1.87	0.41
3:D:1000:THR:HG23	3:D:1041:LEU:HD11	2.02	0.41
2:C:1107:ASN:HA	2:C:1108:PRO:HD3	1.94	0.41
3:D:709:HIS:CD2	3:D:711:LEU:HB2	2.55	0.41
3:D:1100:ASP:O	3:D:1103:HIS:HD2	2.03	0.41
2:C:207:LEU:HD13	2:C:221:LEU:HD21	2.03	0.41
2:C:755:LEU:HB3	2:C:825:VAL:HG21	2.02	0.41
3:D:483:HIS:CE1	3:D:488:ARG:HD3	2.56	0.41
3:D:750:PRO:HG2	3:D:756:GLN:NE2	2.36	0.41
3:D:935:LYS:HE2	3:D:935:LYS:HB3	1.91	0.41
2:C:63:GLY:HA3	2:C:100:LEU:HD21	2.03	0.41
2:C:167:LYS:O	2:C:169:GLY:N	2.52	0.41
2:C:627:ARG:NH2	2:C:640:ARG:HG3	2.35	0.41
3:D:22:SER:HB2	3:D:92:HIS:HB3	2.02	0.41
3:D:106:LYS:HA	3:D:106:LYS:HD2	1.81	0.41
3:D:1336:LEU:HB2	3:D:1344:VAL:HG21	2.01	0.41
2:C:610:ARG:HG2	2:C:611:ILE:N	2.35	0.41
3:D:238:PRO:HD3	3:D:318:ARG:HG3	2.03	0.41
5:F:79:ASP:HA	5:F:80:PRO:HD3	1.95	0.41
8:H:17:DC:H2''	8:H:18:DA:C8	2.56	0.41
2:C:184:MET:HE2	2:C:186:VAL:HG21	2.03	0.41
8:H:19:DC:H6	8:H:19:DC:H2'	1.77	0.41
3:D:490:ALA:O	3:D:494:LYS:HG3	2.21	0.41
3:D:517:VAL:HA	3:D:518:PRO:HD3	1.94	0.41
5:F:330:GLY:HA2	5:F:333:ILE:HD12	2.02	0.41
2:C:713:ARG:CZ	3:D:533:GLY:HA3	2.51	0.41
2:C:892:LEU:HB2	2:C:990:GLY:HA2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:161:LEU:HB3	3:D:452:ILE:CD1	2.51	0.41
3:D:405:ASP:CG	3:D:406:ASP:H	2.29	0.41
3:D:729:HIS:N	11:D:2102:HOH:O	2.53	0.41
4:E:83:ASP:OD1	4:E:83:ASP:N	2.52	0.41
1:B:94:LEU:HD11	1:B:96:THR:O	2.21	0.41
2:C:976:ASP:OD1	2:C:978:ARG:HD3	2.21	0.41
3:D:12:LEU:HA	3:D:12:LEU:HD23	1.89	0.41
3:D:597:ASP:O	3:D:599:PRO:HD3	2.21	0.41
3:D:743:ASP:CG	7:I:7:A:O3'	2.64	0.41
3:D:1350:GLU:OE2	3:D:1357:ARG:NH1	2.46	0.41
4:E:68:LEU:HD12	4:E:68:LEU:HA	1.88	0.41
5:F:408:LEU:HD23	5:F:408:LEU:HA	1.90	0.41
1:A:91:ASN:HA	1:A:92:PRO:HD3	1.88	0.40
1:B:94:LEU:O	1:B:146:ARG:NH2	2.46	0.40
2:C:328:LEU:HA	2:C:328:LEU:HD23	1.76	0.40
2:C:1095:LEU:HD13	3:D:103:TRP:CH2	2.56	0.40
2:C:206:THR:O	2:C:210:GLU:HB2	2.20	0.40
3:D:373:PRO:HA	3:D:376:GLU:HG3	2.03	0.40
3:D:1018:ASN:O	3:D:1022:VAL:HG23	2.21	0.40
1:B:101:LEU:HD11	1:B:113:ASP:HB2	2.03	0.40
2:C:6:PHE:CD1	2:C:909:ALA:HB2	2.57	0.40
2:C:838:LYS:HE2	2:C:997:LEU:HD12	2.04	0.40
2:C:853:LEU:HB2	2:C:858:MET:SD	2.61	0.40
3:D:134:VAL:HG23	3:D:151:GLN:O	2.21	0.40
3:D:372:ASP:HA	3:D:373:PRO:HD3	1.92	0.40
3:D:633:VAL:HB	3:D:740:PHE:CE2	2.55	0.40
3:D:1053:PHE:CE2	3:D:1072:ILE:HG23	2.56	0.40
3:D:1407:LEU:O	3:D:1412:LYS:N	2.53	0.40
5:F:93:LEU:HA	5:F:93:LEU:HD23	1.89	0.40
5:F:295:MET:HB3	5:F:299:TRP:CG	2.55	0.40
8:H:13:DC:C1'	8:H:14:DT:H72	2.52	0.40
1:B:110:LYS:HD2	1:B:126:ASP:O	2.21	0.40
3:D:864:VAL:HG22	3:D:865:THR:H	1.86	0.40
3:D:895:VAL:O	3:D:899:LEU:HG	2.22	0.40
4:E:59:ASN:O	4:E:63:TRP:HD1	2.04	0.40
5:F:338:LEU:HA	5:F:339:PRO:HD3	1.86	0.40
2:C:570:PRO:HB3	2:C:660:ALA:HB2	2.04	0.40
2:C:886:LEU:HD21	3:D:951:ILE:HG12	2.04	0.40
3:D:355:VAL:HG11	3:D:385:VAL:HG21	2.02	0.40
3:D:1057:VAL:HG13	3:D:1069:GLU:CD	2.47	0.40
5:F:361:LEU:HB3	5:F:365:GLU:HG3	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	230/315 (73%)	227 (99%)	3 (1%)	0	100	100
1	B	221/315 (70%)	216 (98%)	4 (2%)	1 (0%)	24	55
2	C	1111/1119 (99%)	1076 (97%)	35 (3%)	0	100	100
3	D	1484/1524 (97%)	1433 (97%)	51 (3%)	0	100	100
4	E	92/99 (93%)	89 (97%)	3 (3%)	0	100	100
5	F	331/443 (75%)	327 (99%)	4 (1%)	0	100	100
All	All	3469/3815 (91%)	3368 (97%)	100 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	8	ALA

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	201/273 (74%)	199 (99%)	2 (1%)	68	76
1	B	196/273 (72%)	191 (97%)	5 (3%)	40	64
2	C	939/941 (100%)	916 (98%)	23 (2%)	43	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	D	1255/1279 (98%)	1217 (97%)	38 (3%)	36	61
4	E	82/88 (93%)	80 (98%)	2 (2%)	43	65
5	F	291/388 (75%)	282 (97%)	9 (3%)	35	60
All	All	2964/3242 (91%)	2885 (97%)	79 (3%)	40	63

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	134	GLU
1	B	10	VAL
1	B	133	GLU
1	B	154	GLU
1	B	188	GLN
1	B	191	ASP
2	C	11	GLU
2	C	27	ARG
2	C	177	GLU
2	C	182	VAL
2	C	217	LEU
2	C	276	LYS
2	C	284	ARG
2	C	348	LEU
2	C	384	GLU
2	C	388	ARG
2	C	390	GLN
2	C	419	THR
2	C	427	VAL
2	C	443	THR
2	C	513	VAL
2	C	595	LEU
2	C	617	ASP
2	C	723	THR
2	C	775	ARG
2	C	807[A]	ARG
2	C	807[B]	ARG
2	C	888	THR
2	C	929	ARG
3	D	40	GLU
3	D	71	LYS
3	D	145	VAL

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Mol	Chain	Res	Type
3	D	191	LEU
3	D	231	VAL
3	D	270	LEU
3	D	311	LEU
3	D	331	VAL
3	D	354	VAL
3	D	443	VAL
3	D	452	ILE
3	D	471	GLU
3	D	632	VAL
3	D	633	VAL
3	D	650	LEU
3	D	666	ILE
3	D	687	VAL
3	D	741	ASP
3	D	747	VAL
3	D	754	PHE
3	D	778	LEU
3	D	810	GLU
3	D	894	LYS
3	D	907	GLU
3	D	921	ARG
3	D	922	LEU
3	D	940	THR
3	D	986	ARG
3	D	1129	THR
3	D	1188	VAL
3	D	1207	TYR
3	D	1231	GLU
3	D	1234	THR
3	D	1252	ILE
3	D	1283	ILE
3	D	1286	THR
3	D	1295	GLU
3	D	1501	GLU
4	E	49	GLN
4	E	50	THR
5	F	88	ILE
5	F	122	LEU
5	F	193	ARG
5	F	201	LYS
5	F	218	GLN

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Mol	Chain	Res	Type
5	F	295	MET
5	F	369	LEU
5	F	373	LYS
5	F	380	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	HIS
1	B	156	HIS
1	B	212	ASN
2	C	99	GLN
2	C	343	GLN
2	C	390	GLN
2	C	543	ASN
2	C	834	GLN
2	C	1026	GLN
2	C	1047	HIS
3	D	66	GLN
3	D	442	ASN
3	D	507	ASN
3	D	669	ASN
3	D	717	GLN
3	D	724	GLN
3	D	762	GLN
3	D	816	HIS
3	D	991	GLN
3	D	1046	GLN
3	D	1075	HIS
3	D	1103	HIS
3	D	1124	GLN
3	D	1172	HIS
3	D	1359	GLN
3	D	1441	GLN
3	D	1445	HIS
5	F	83	GLN
5	F	86	HIS
5	F	218	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	I	6/7 (85%)	1 (16%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
7	I	2	C

There are no RNA pucker outliers to report.

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 6 ligands modelled in this entry, 6 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	231/315 (73%)	0.02	4 (1%) 69 50	74, 114, 142, 189	1 (0%)
1	B	223/315 (70%)	-0.24	1 (0%) 88 79	66, 99, 137, 157	0
2	C	1112/1119 (99%)	0.08	20 (1%) 67 49	55, 114, 176, 217	3 (0%)
3	D	1486/1524 (97%)	-0.03	21 (1%) 73 55	50, 97, 174, 205	3 (0%)
4	E	94/99 (94%)	0.27	2 (2%) 63 44	76, 124, 176, 182	0
5	F	335/443 (75%)	0.29	8 (2%) 59 40	88, 138, 224, 236	0
6	G	18/21 (85%)	-0.35	0 100 100	73, 104, 196, 202	0
7	I	7/7 (100%)	-0.72	0 100 100	69, 76, 122, 132	0
8	H	25/28 (89%)	0.27	0 100 100	102, 124, 181, 209	0
All	All	3531/3871 (91%)	0.03	56 (1%) 70 52	50, 110, 179, 236	7 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	F	196	VAL	4.4
3	D	1073	SER	4.2
3	D	282	TYR	4.1
3	D	353	VAL	3.8
3	D	409	VAL	3.7
5	F	192	LEU	3.6
3	D	368	VAL	3.4
2	C	47	ALA	3.2
2	C	68	PHE	3.0
3	D	225	LEU	3.0
3	D	355	VAL	3.0
5	F	191	ASN	3.0
3	D	211	VAL	2.9
2	C	65	VAL	2.9
3	D	269	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
2	C	66	LEU	2.9
2	C	43	GLY	2.8
2	C	1091	GLU	2.8
2	C	998	TYR	2.8
2	C	98	LEU	2.8
2	C	587	VAL	2.7
2	C	372	LEU	2.6
5	F	190	ALA	2.6
3	D	520	LEU	2.6
5	F	331	ASP	2.6
2	C	46	ALA	2.6
3	D	367	ILE	2.6
5	F	239	ALA	2.5
5	F	217	ASN	2.5
2	C	138	SER	2.5
3	D	1495	ILE	2.5
2	C	103	LYS	2.4
3	D	304	LEU	2.4
1	A	79	ILE	2.4
2	C	536	PRO	2.4
1	A	129	ILE	2.3
2	C	467	ILE	2.3
3	D	305	ALA	2.3
3	D	324	ALA	2.3
5	F	199	ALA	2.3
2	C	101	ILE	2.2
2	C	373	VAL	2.2
2	C	25	SER	2.2
2	C	64	LEU	2.2
4	E	32	ARG	2.2
3	D	647	ARG	2.2
1	A	73	GLU	2.1
2	C	254	VAL	2.1
3	D	184	GLU	2.1
3	D	207	PHE	2.1
1	A	131	THR	2.1
1	B	228	PRO	2.1
3	D	322	VAL	2.0
4	E	51	LEU	2.0
3	D	290	PRO	2.0
3	D	1283	ILE	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
9	MG	B	2001	1/1	0.62	0.09	118,118,118,118	0
9	MG	D	2004	1/1	0.81	0.25	65,65,65,65	0
9	MG	D	2003	1/1	0.95	0.07	53,53,53,53	0
10	ZN	D	2002	1/1	0.96	0.09	163,163,163,163	0
9	MG	F	2001	1/1	0.98	0.04	135,135,135,135	0
10	ZN	D	2001	1/1	0.99	0.02	73,73,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.