



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:17 PM UTC

PDB ID : 6KQE / pdb_00006kqe
Title : Thermus thermophilus initial transcription complex comprising sigma A and 5'-OH RNA of 4 nt
Authors : Zhang, Y.; Li, L.; Ebright, R.H.
Deposited on : 2019-08-17
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

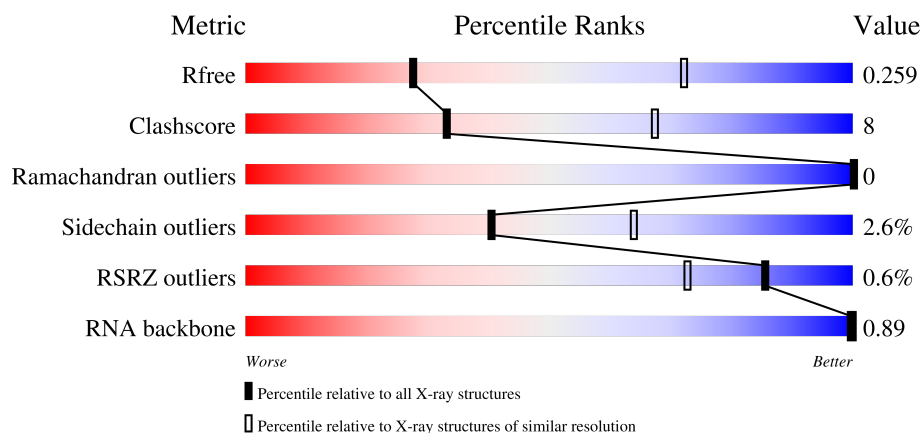
1 Overall quality at a glance

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	
1	B	315	
2	C	1119	
3	D	1524	

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Mol	Chain	Length	Quality of chain
4	E	99	79%16%5%
5	F	443	% 65%13%22%
6	G	21	48%33%19%
7	H	27	33%56%11%
8	I	4	100%

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 28688 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	0	0
			1809	1155	315	337	2			
1	B	227	Total	C	N	O	S	0	0	0
			1784	1140	309	333	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	0	0
			8762	5544	1559	1635	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
			759	484	132	139	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	346	Total	C	N	O	S	0	0	0
			2801	1767	506	524	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	17	Total	C	N	O	P	0	0	0
			350	166	68	100	16			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	24	Total	C	N	O	P	0	0	0
			495	236	94	142	23			

- Molecule 8 is a RNA chain called RNA (5'-R(*UP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	I	4	Total	C	N	O	P	0	0	0
			82	38	15	26	3			

- 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

- 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	8	Total O 8 8	0	0
11	B	4	Total O 4 4	0	0
11	C	23	Total O 23 23	0	0
11	D	39	Total O 39 39	0	0
11	E	2	Total O 2 2	0	0
11	F	9	Total O 9 9	0	0
11	G	3	Total O 3 3	0	0
11	H	1	Total O 1 1	0	0
11	I	1	Total O 1 1	0	0

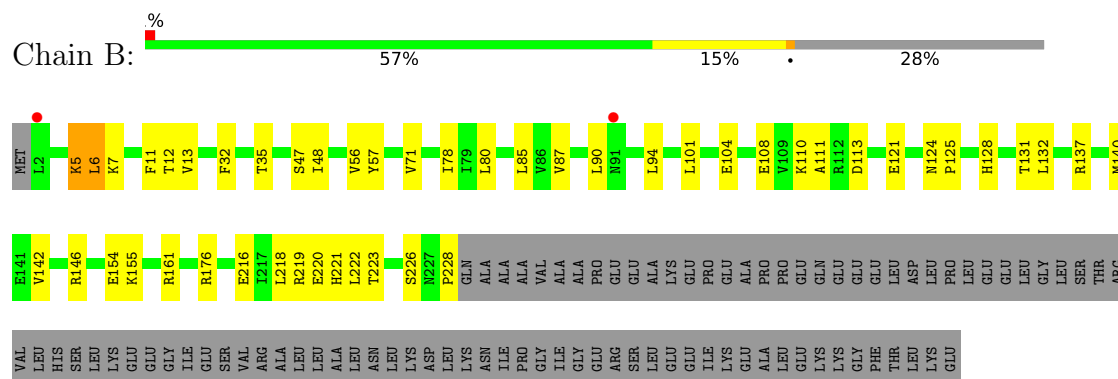
3 Residue-property plots [i](#)

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.

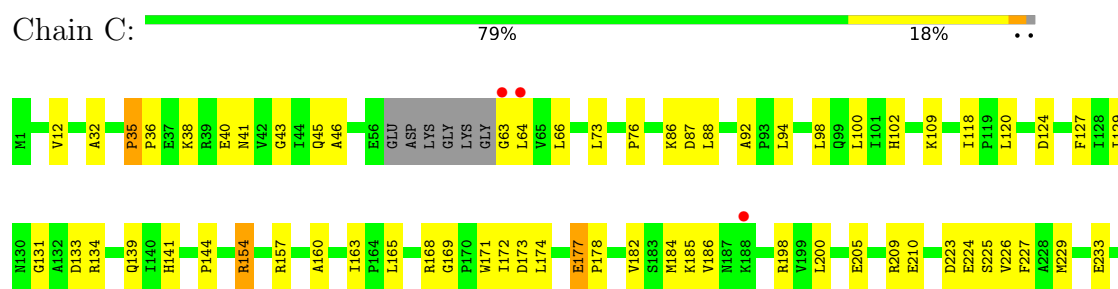
- Molecule 1: DNA-directed RNA polymerase subunit alpha

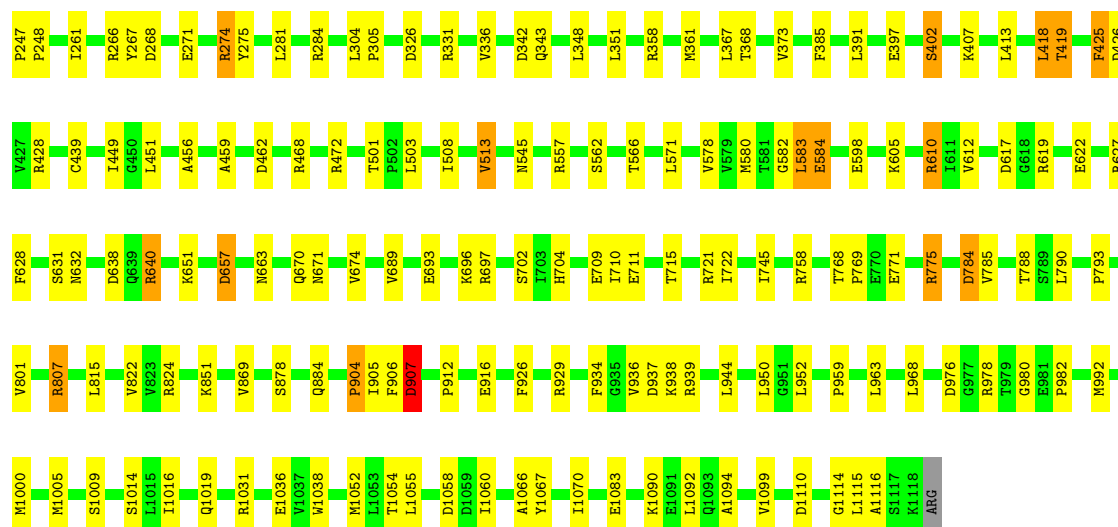


- Molecule 1: DNA-directed RNA polymerase subunit alpha

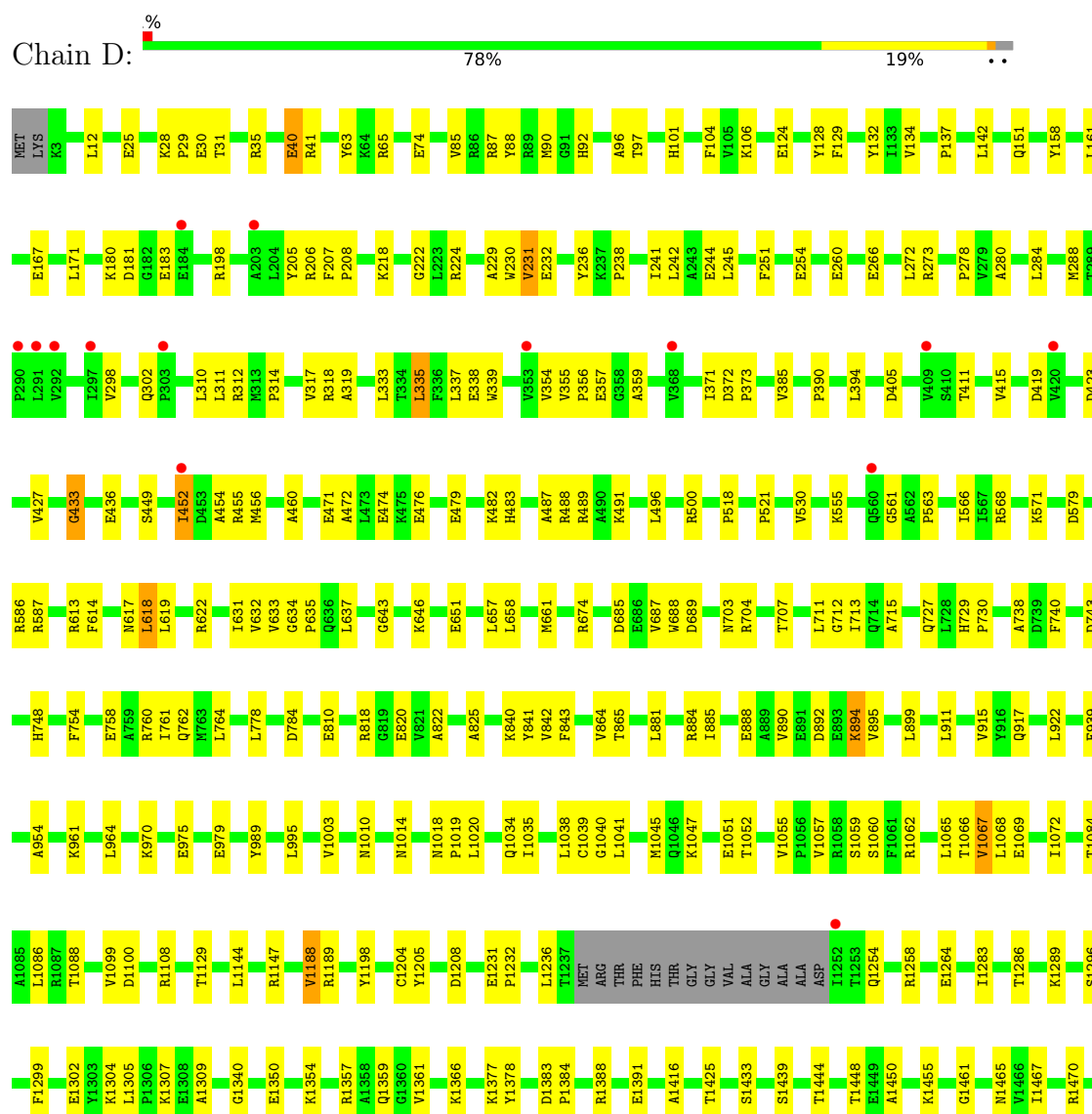


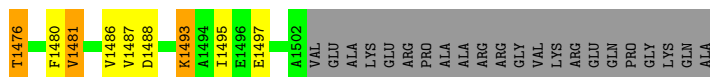
- Molecule 2: DNA-directed RNA polymerase subunit beta



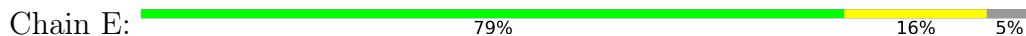


• Molecule 3: DNA-directed RNA polymerase subunit beta'

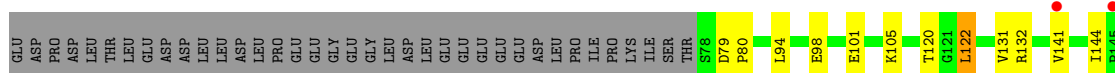
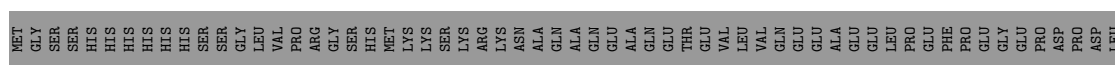




- Molecule 4: DNA-directed RNA polymerase subunit omega



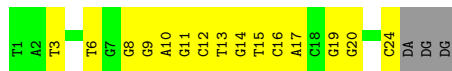
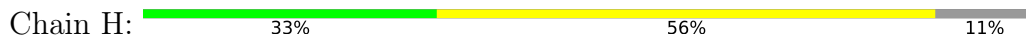
- Molecule 5: RNA polymerase sigma factor SigA



- 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: DNA (5'-D(*TP*AP*TP*AP*AP*TP*GP*GP*GP*AP*GP*CP*TP*GP*TP*CP*AP*CP*GP*GP*AP*TP*GP*CP*AP*GP*G*)-3')



- Molecule 8: RNA (5'-R(*UP*CP*GP*A)-3')



There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	184.57Å 103.76Å 295.98Å 90.00° 98.70° 90.00°	Depositor
Resolution (Å)	48.95 – 3.30 48.95 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.95-3.30) 99.9 (48.95-3.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.217 , 0.258 0.217 , 0.259	Depositor DCC
R_{free} test set	4156 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	85.9	Xtriage
Anisotropy	0.523	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 40.4	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	0.94	EDS
Total number of atoms	28688	wwPDB-VP
Average B, all atoms (Å ²)	95.0	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: ZN, MG

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	0.31	0/1841	0.86	6/2504 (0.2%)
1	B	0.32	0/1816	0.82	4/2469 (0.2%)
2	C	0.36	0/8929	0.80	15/12078 (0.1%)
3	D	0.32	0/11963	0.79	10/16173 (0.1%)
4	E	0.30	0/773	0.76	0/1042
5	F	0.30	0/2846	0.76	2/3830 (0.1%)
6	G	0.16	0/393	0.46	0/606
7	H	0.21	0/556	0.61	0/858
8	I	0.13	0/91	0.26	0/140
All	All	0.33	0/29208	0.79	37/39700 (0.1%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1231	GLU	CA-C-N	-8.88	112.68	121.65
3	D	1231	GLU	C-N-CA	-8.88	112.68	121.65
1	A	172	SER	CA-C-N	7.51	127.69	119.87
1	A	172	SER	C-N-CA	7.51	127.69	119.87
3	D	1340	GLY	CA-C-N	7.31	127.66	119.32
3	D	1340	GLY	C-N-CA	7.31	127.66	119.32
3	D	433	GLY	N-CA-C	7.17	120.23	110.69
1	A	8	ALA	CA-C-N	6.35	126.87	120.14
1	A	8	ALA	C-N-CA	6.35	126.87	120.14
1	A	115	LEU	CA-C-N	5.88	125.81	119.76
1	A	115	LEU	C-N-CA	5.88	125.81	119.76
2	C	907	ASP	N-CA-C	-5.79	98.47	110.80
3	D	1481	VAL	N-CA-C	5.51	118.28	112.83
2	C	35	PRO	CA-C-N	5.49	125.84	119.47
2	C	35	PRO	C-N-CA	5.49	125.84	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	419	THR	N-CA-C	-5.45	102.42	110.48
1	B	48	ILE	CA-C-N	5.42	125.69	119.83
1	B	48	ILE	C-N-CA	5.42	125.69	119.83
3	D	1208	ASP	N-CA-C	-5.36	99.78	108.41
1	B	47	SER	N-CA-C	5.31	119.86	113.17
2	C	154	ARG	CA-C-N	5.29	124.48	118.97
2	C	154	ARG	C-N-CA	5.29	124.48	118.97
2	C	980	GLY	N-CA-C	-5.26	107.28	114.64
3	D	617	ASN	N-CA-C	5.20	119.72	113.17
2	C	904	PRO	CA-C-O	-5.18	115.69	121.23
2	C	793	PRO	O-C-N	5.17	123.69	121.31
2	C	418	LEU	N-CA-C	5.11	117.64	110.23
2	C	271	GLU	N-CA-C	-5.10	105.80	111.36
2	C	439	CYS	CA-C-N	5.09	124.70	119.56
2	C	439	CYS	C-N-CA	5.09	124.70	119.56
2	C	169	GLY	CA-C-N	5.07	125.51	120.14
2	C	169	GLY	C-N-CA	5.07	125.51	120.14
3	D	740	PHE	N-CA-C	5.06	119.14	112.92
5	F	296	GLY	CA-C-N	5.04	124.94	119.85
5	F	296	GLY	C-N-CA	5.04	124.94	119.85
1	B	226	SER	N-CA-C	5.01	116.60	111.03
3	D	740	PHE	CB-CA-C	5.01	119.03	110.56

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	1809	0	1863	28	0
1	B	1784	0	1830	39	0
2	C	8762	0	8855	144	0
3	D	11751	0	11993	191	0
4	E	759	0	771	12	0
5	F	2801	0	2871	35	0
6	G	350	0	192	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	H	495	0	272	41	0
8	I	82	0	44	0	0
9	B	1	0	0	0	0
9	D	2	0	0	0	0
10	D	2	0	0	0	0
11	A	8	0	0	1	0
11	B	4	0	0	0	0
11	C	23	0	0	1	0
11	D	39	0	0	1	0
11	E	2	0	0	1	0
11	F	9	0	0	0	0
11	G	3	0	0	0	0
11	H	1	0	0	0	0
11	I	1	0	0	0	0
All	All	28688	0	28691	447	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (447) 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)
2:C:425:PHE:CE1	3:D:1086:LEU:HD12	1.74	1.23
7:H:16:DC:H2''	7:H:17:DA:C8	1.78	1.18
7:H:19:DG:H2''	7:H:20:DG:C8	1.80	1.16
6:G:4:DG:H2''	6:G:5:DC:C6	1.82	1.13
6:G:4:DG:H2''	6:G:5:DC:C5	1.82	1.13
7:H:11:DG:H3'	7:H:12:DC:H5''	1.26	1.10
2:C:425:PHE:HE1	3:D:1086:LEU:CD1	1.65	1.08
2:C:425:PHE:HE1	3:D:1086:LEU:HD12	0.98	1.07
2:C:906:PHE:CD2	3:D:1067:VAL:HG12	1.95	1.01
7:H:19:DG:C5	7:H:20:DG:C6	2.52	0.97
6:G:11:DT:H2''	6:G:12:DG:C8	2.00	0.97
6:G:4:DG:C2'	6:G:5:DC:C5	2.47	0.97
7:H:19:DG:H2''	7:H:20:DG:H8	1.29	0.90
1:B:7:LYS:HE2	1:B:7:LYS:N	1.88	0.88
7:H:19:DG:C2'	7:H:20:DG:C8	2.56	0.88
7:H:11:DG:H3'	7:H:12:DC:C5'	2.04	0.87
2:C:418:LEU:HD21	7:H:14:DG:C8	2.10	0.87
7:H:16:DC:C2'	7:H:17:DA:C8	2.59	0.86
7:H:15:DT:C4	7:H:16:DC:N4	2.44	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:628:PHE:H	2:C:638:ASP:HB3	1.42	0.84
7:H:16:DC:H2''	7:H:17:DA:N7	1.95	0.79
2:C:418:LEU:HD21	7:H:14:DG:N7	1.97	0.78
1:B:5:LYS:HD2	1:B:5:LYS:O	1.85	0.77
2:C:906:PHE:CE2	3:D:1067:VAL:HA	2.19	0.77
5:F:338:LEU:HD23	5:F:339:PRO:HD2	1.67	0.76
2:C:906:PHE:HE2	3:D:1067:VAL:N	1.83	0.76
1:A:228:PRO:HA	1:B:11:PHE:CD2	2.20	0.76
2:C:428:ARG:HD3	2:C:449:ILE:O	1.86	0.75
7:H:8:DG:H2''	7:H:9:DG:O5'	1.87	0.75
6:G:4:DG:H1	7:H:24:DC:H42	1.35	0.74
2:C:139:GLN:HB2	2:C:391:LEU:HD21	1.70	0.73
2:C:906:PHE:CE2	3:D:1067:VAL:CA	2.73	0.72
2:C:425:PHE:CE1	3:D:1086:LEU:CD1	2.53	0.71
1:B:6:LEU:C	1:B:7:LYS:HE2	2.15	0.71
7:H:19:DG:C6	7:H:20:DG:C6	2.79	0.71
2:C:906:PHE:CE2	3:D:1067:VAL:N	2.59	0.70
2:C:904:PRO:HB2	2:C:907:ASP:O	1.92	0.70
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.74	0.70
3:D:134:VAL:HG22	3:D:151:GLN:H	1.56	0.70
1:B:6:LEU:HD13	1:B:6:LEU:O	1.93	0.69
2:C:266:ARG:NH1	7:H:11:DG:O6	2.26	0.69
6:G:4:DG:C2'	6:G:5:DC:H5	2.06	0.69
2:C:226:VAL:HA	2:C:229:MET:HE2	1.73	0.68
3:D:284:LEU:HD22	3:D:288:MET:HE2	1.75	0.68
1:B:6:LEU:O	1:B:6:LEU:HD22	1.93	0.68
5:F:316:SER:HB3	5:F:319:THR:HG23	1.75	0.68
7:H:15:DT:C4	7:H:16:DC:C4	2.82	0.67
5:F:193:ARG:HB2	7:H:6:DT:H1'	1.77	0.67
3:D:208:PRO:HA	3:D:390:PRO:HA	1.78	0.66
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.78	0.66
1:B:5:LYS:HB2	1:B:7:LYS:HZ1	1.62	0.65
3:D:224:ARG:H	3:D:251:PHE:HE1	1.44	0.65
5:F:361:LEU:HB3	5:F:365:GLU:HG3	1.79	0.65
7:H:19:DG:C4	7:H:20:DG:C5	2.85	0.65
3:D:132:TYR:OH	3:D:568:ARG:NH1	2.30	0.65
2:C:413:LEU:HD21	2:C:451:LEU:HD13	1.79	0.64
3:D:65:ARG:NH1	5:F:378:GLY:O	2.31	0.64
2:C:906:PHE:HD2	3:D:1067:VAL:HG12	1.59	0.64
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.80	0.64
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:612:VAL:HG22	2:C:622:GLU:HG3	1.80	0.64
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.79	0.63
3:D:1236:LEU:HA	3:D:1359:GLN:HG3	1.79	0.63
2:C:711:GLU:HG2	2:C:822:VAL:HG22	1.79	0.63
2:C:906:PHE:HE2	3:D:1066:THR:C	2.07	0.63
2:C:904:PRO:HD2	2:C:907:ASP:O	1.99	0.63
7:H:15:DT:N3	7:H:16:DC:C4	2.67	0.63
3:D:1040:GLY:O	3:D:1060:SER:HB3	1.99	0.63
3:D:619:LEU:HD11	3:D:1439:SER:HB3	1.79	0.63
1:A:36:LEU:HD11	1:B:221:HIS:HB3	1.80	0.62
6:G:4:DG:H2'	6:G:5:DC:C5	2.33	0.62
2:C:168:ARG:HD3	2:C:268:ASP:HB3	1.80	0.62
2:C:950:LEU:HB3	2:C:952:LEU:HD13	1.81	0.62
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.80	0.62
3:D:657:LEU:HG	3:D:661:MET:HE2	1.82	0.62
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.33	0.62
2:C:1067:TYR:OH	3:D:674[A]:ARG:NH1	2.32	0.62
3:D:685:ASP:HA	3:D:688:TRP:HD1	1.64	0.62
3:D:1439:SER:OG	3:D:1467:ILE:HD11	2.00	0.62
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.81	0.62
2:C:12:VAL:HG21	2:C:472:ARG:HD3	1.82	0.62
7:H:19:DG:C6	7:H:20:DG:N1	2.69	0.61
5:F:365:GLU:HB2	5:F:404:ALA:HB2	1.82	0.61
1:B:80:LEU:HD21	3:D:842:VAL:HG12	1.81	0.61
3:D:41:ARG:NH2	11:D:2104:HOH:O	2.32	0.61
2:C:397:GLU:HG3	2:C:631:SER:HB2	1.82	0.61
3:D:954:ALA:O	3:D:1062:ARG:NH2	2.34	0.60
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.83	0.60
5:F:397:ILE:HD12	5:F:400:ILE:HD11	1.82	0.60
5:F:400:ILE:HA	5:F:403:LYS:HG2	1.83	0.59
6:G:11:DT:C2'	6:G:12:DG:C8	2.82	0.59
3:D:1465:ASN:OD1	3:D:1470:ARG:NH1	2.35	0.59
2:C:1083:GLU:OE2	3:D:87:ARG:NH2	2.35	0.59
3:D:622:ARG:NH1	6:G:17:DG:OP1	2.36	0.59
6:G:4:DG:H2'	6:G:5:DC:H5	1.66	0.58
2:C:689:VAL:HG13	2:C:851:LYS:HB3	1.83	0.58
3:D:272:LEU:HB2	3:D:280:ALA:HB3	1.85	0.58
3:D:711:LEU:HD13	3:D:778:LEU:HD23	1.84	0.58
1:B:220:GLU:O	1:B:223:THR:OG1	2.20	0.58
3:D:489:ARG:NH1	3:D:1391:GLU:OE2	2.35	0.57
3:D:433:GLY:HA2	3:D:449:SER:H	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:707:THR:HG23	3:D:712:GLY:HA3	1.85	0.57
7:H:19:DG:C5	7:H:20:DG:C5	2.93	0.57
3:D:134:VAL:HG12	3:D:454:ALA:HB2	1.87	0.57
3:D:167:GLU:OE2	3:D:198:ARG:NH1	2.38	0.57
1:B:216:GLU:OE1	1:B:219:ARG:NH2	2.28	0.56
2:C:939:ARG:HG2	2:C:982:PRO:HD3	1.87	0.56
2:C:628:PHE:H	2:C:638:ASP:CB	2.17	0.56
3:D:273:ARG:HB3	3:D:278:PRO:HA	1.87	0.56
1:B:176:ARG:NH2	3:D:888:GLU:OE1	2.36	0.56
2:C:425:PHE:HD1	2:C:425:PHE:N	2.03	0.56
2:C:1092:LEU:HD13	2:C:1099:VAL:HG21	1.87	0.56
2:C:425:PHE:N	2:C:425:PHE:CD1	2.73	0.56
2:C:129:ILE:HB	2:C:134:ARG:HD2	1.88	0.56
2:C:402:SER:HA	2:C:566:THR:HG23	1.87	0.56
2:C:905:ILE:HG13	2:C:905:ILE:O	2.04	0.56
7:H:12:DC:H2''	7:H:13:DT:C6	2.41	0.56
7:H:17:DA:OP2	7:H:17:DA:H2'	2.06	0.56
7:H:19:DG:N7	7:H:20:DG:O6	2.39	0.56
1:B:6:LEU:C	1:B:6:LEU:HD22	2.31	0.55
2:C:223:ASP:OD1	2:C:225:SER:OG	2.23	0.55
3:D:171:LEU:HD12	3:D:390:PRO:HG2	1.88	0.55
3:D:864:VAL:HG12	3:D:865:THR:H	1.71	0.55
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.88	0.55
2:C:35:PRO:HG2	2:C:38:LYS:HD2	1.89	0.55
3:D:890:VAL:HB	3:D:922:LEU:HD13	1.89	0.55
2:C:41:ASN:O	2:C:46:ALA:HB2	2.07	0.55
2:C:165:LEU:HB2	2:C:168:ARG:HG3	1.87	0.55
2:C:768:THR:OG1	2:C:771:GLU:OE1	2.25	0.54
2:C:160:ALA:HB3	2:C:174:LEU:HB2	1.88	0.54
2:C:501:THR:HG21	2:C:513:VAL:HG23	1.90	0.54
1:A:183:ASP:HA	2:C:938:LYS:HE3	1.89	0.54
2:C:1016:ILE:O	3:D:87:ARG:NH1	2.41	0.54
2:C:545:ASN:HB3	2:C:583:LEU:HD22	1.89	0.53
1:B:6:LEU:H	1:B:6:LEU:CD1	2.21	0.53
2:C:1067:TYR:OH	3:D:674[B]:ARG:NH1	2.42	0.53
3:D:890:VAL:HG23	3:D:892:ASP:H	1.73	0.53
2:C:758:ARG:HH21	2:C:788:THR:HB	1.73	0.53
2:C:944:LEU:HD21	2:C:963:LEU:HD23	1.91	0.53
3:D:703:ASN:HB2	3:D:713:ILE:HG12	1.91	0.53
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.90	0.53
3:D:1254:GLN:HB3	3:D:1258:ARG:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:709:GLU:OE2	2:C:824:ARG:NH1	2.42	0.53
2:C:172:ILE:HG12	2:C:186:VAL:HG22	1.90	0.53
3:D:137:PRO:HA	3:D:452:ILE:HG13	1.89	0.53
3:D:843:PHE:HE2	3:D:864:VAL:HG11	1.73	0.52
5:F:80:PRO:HB2	5:F:210:LEU:HD11	1.91	0.52
2:C:210:GLU:HG2	2:C:304:LEU:HD21	1.90	0.52
2:C:598:GLU:O	2:C:651:LYS:NZ	2.36	0.52
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.91	0.52
2:C:674:VAL:HG23	2:C:869:VAL:HB	1.91	0.52
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.92	0.52
4:E:37:ASN:HD22	4:E:89:MET:HE1	1.74	0.52
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.90	0.52
2:C:1052:MET:HE1	3:D:748:HIS:HB3	1.92	0.52
3:D:411:THR:HG23	3:D:436:GLU:HA	1.92	0.52
6:G:12:DG:H5"	6:G:12:DG:H8	1.75	0.52
2:C:168:ARG:O	2:C:267:TYR:HA	2.10	0.52
3:D:314:PRO:HB2	3:D:317:VAL:HG12	1.91	0.52
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.45	0.51
1:B:94:LEU:O	1:B:146:ARG:NH2	2.43	0.51
3:D:218:LYS:HG2	3:D:338:GLU:HG2	1.91	0.51
3:D:563:PRO:HD2	3:D:566:ILE:HD12	1.91	0.51
2:C:173:ASP:HB2	2:C:185:LYS:HB3	1.91	0.51
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.92	0.51
3:D:843:PHE:CE2	3:D:864:VAL:HG11	2.45	0.51
7:H:16:DC:H2"	7:H:17:DA:H8	1.59	0.51
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.93	0.51
1:B:104:GLU:OE2	1:B:137:ARG:NH1	2.44	0.51
3:D:30:GLU:OE1	3:D:40:GLU:HG2	2.11	0.51
2:C:571:LEU:HD23	2:C:702:SER:HB3	1.93	0.51
3:D:1144:LEU:O	3:D:1147:ARG:HG3	2.11	0.51
5:F:237:THR:HG21	7:H:3:DT:H5"	1.91	0.51
5:F:276:ARG:O	5:F:279:GLN:HG3	2.11	0.51
7:H:16:DC:C2'	7:H:17:DA:N7	2.69	0.51
3:D:405:ASP:HB3	3:D:423:ASP:HA	1.92	0.51
1:A:209:GLU:O	1:A:213:GLN:HG2	2.10	0.50
2:C:557:ARG:NH2	11:C:1202:HOH:O	2.38	0.50
3:D:455:ARG:HB2	3:D:460:ALA:HB2	1.92	0.50
1:B:110:LYS:HD3	1:B:128:HIS:HA	1.94	0.50
3:D:1189:ARG:HB3	3:D:1204:CYS:HA	1.93	0.50
3:D:658:LEU:HA	3:D:661:MET:HE3	1.94	0.50
3:D:840:LYS:HE3	3:D:841:TYR:CZ	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:H:16:DC:H1'	7:H:17:DA:C8	2.47	0.50
7:H:19:DG:C5	7:H:20:DG:O6	2.64	0.50
1:B:5:LYS:HB2	1:B:7:LYS:NZ	2.27	0.50
5:F:395:GLU:OE2	5:F:398:ARG:NH2	2.44	0.50
3:D:643:GLY:HA3	3:D:727:GLN:HB2	1.94	0.50
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.12	0.50
7:H:19:DG:N7	7:H:20:DG:C6	2.79	0.50
2:C:1031:ARG:HG2	6:G:16:DC:H5''	1.93	0.49
3:D:241:ILE:HA	3:D:312:ARG:HB3	1.93	0.49
3:D:894:LYS:H	3:D:894:LYS:HD3	1.77	0.49
5:F:163:LEU:HD13	5:F:174:LEU:HD13	1.94	0.49
2:C:63:GLY:HA3	2:C:100:LEU:HD21	1.94	0.49
2:C:710:ILE:HD12	2:C:790:LEU:HB2	1.93	0.49
2:C:937:ASP:OD1	2:C:939:ARG:HD3	2.12	0.49
3:D:180:LYS:NZ	3:D:357:GLU:OE1	2.34	0.49
3:D:566:ILE:HD11	5:F:192:LEU:HD21	1.94	0.49
1:A:228:PRO:HG3	1:B:11:PHE:HE2	1.78	0.49
5:F:101:GLU:HG2	5:F:105:LYS:HE2	1.95	0.49
3:D:97:THR:HG21	3:D:571:LYS:HG2	1.94	0.49
1:A:121:GLU:OE1	11:A:401:HOH:O	2.20	0.49
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.93	0.49
4:E:44:GLU:OE2	4:E:72:ARG:NH1	2.44	0.49
3:D:244:GLU:HG3	3:D:310:LEU:HG	1.95	0.48
2:C:1031:ARG:NE	6:G:16:DC:OP1	2.43	0.48
2:C:1036:GLU:OE2	2:C:1036:GLU:N	2.47	0.48
3:D:224:ARG:NE	3:D:254:GLU:OE2	2.25	0.48
1:A:31:GLY:N	1:A:193:ASP:OD2	2.46	0.48
2:C:351:LEU:HD11	2:C:373:VAL:HG13	1.95	0.48
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.95	0.48
7:H:16:DC:C1'	7:H:17:DA:C8	2.96	0.48
3:D:1488:ASP:OD1	3:D:1488:ASP:N	2.45	0.48
3:D:101:HIS:HB3	3:D:104:PHE:HD2	1.77	0.48
1:B:111:ALA:HB3	1:B:125:PRO:HA	1.96	0.48
3:D:129:PHE:CD2	3:D:456:MET:HB3	2.49	0.48
3:D:1383:ASP:HB3	3:D:1416:ALA:HB3	1.96	0.48
2:C:326:ASP:HA	2:C:331:ARG:HD2	1.95	0.48
3:D:864:VAL:HG12	3:D:865:THR:N	2.29	0.48
3:D:1232:PRO:HG3	3:D:1361:VAL:HG11	1.96	0.48
3:D:487:ALA:O	3:D:491:LYS:HG2	2.13	0.48
1:B:57:TYR:CG	1:B:161:ARG:HD2	2.48	0.48
2:C:198:ARG:HE	2:C:227:PHE:HA	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:397:GLU:HG2	2:C:632:ASN:HB2	1.95	0.47
2:C:580:MET:SD	2:C:584:GLU:HG3	2.54	0.47
1:A:64:GLU:HG3	1:A:79:ILE:HD12	1.95	0.47
3:D:970:LYS:HD3	3:D:995:LEU:HD13	1.96	0.47
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.96	0.47
7:H:19:DG:C8	7:H:20:DG:N7	2.82	0.47
1:A:57:TYR:CG	1:A:161:ARG:HD2	2.49	0.47
1:B:78:ILE:HG21	1:B:140:MET:HE1	1.97	0.47
3:D:500:ARG:NH1	3:D:1388:ARG:O	2.48	0.47
1:A:228:PRO:CA	1:B:11:PHE:CD2	2.95	0.47
2:C:426:ASP:OD1	2:C:426:ASP:N	2.48	0.47
2:C:627:ARG:HA	2:C:638:ASP:HB2	1.97	0.47
3:D:181:ASP:HB2	3:D:205:TYR:CD2	2.50	0.47
4:E:40:LEU:HG	4:E:67:GLU:HG2	1.96	0.47
5:F:79:ASP:OD2	7:H:8:DG:N1	2.35	0.47
5:F:122:LEU:HD21	5:F:159:ILE:HD12	1.97	0.46
5:F:328:PHE:HB2	5:F:331:ASP:OD2	2.14	0.46
3:D:975:GLU:O	3:D:979:GLU:HG2	2.15	0.46
2:C:503:LEU:HD23	2:C:508:ILE:HA	1.97	0.46
1:B:32:PHE:HA	1:B:35:THR:HB	1.98	0.46
2:C:32:ALA:HB2	2:C:73:LEU:HD12	1.97	0.46
2:C:154:ARG:HE	2:C:157:ARG:HG3	1.80	0.46
3:D:704:ARG:HD2	3:D:738:ALA:HB2	1.98	0.46
1:A:6:LEU:HD11	1:A:27:PRO:HG2	1.98	0.46
1:B:85:LEU:HG	1:B:87:VAL:HG23	1.97	0.46
1:B:155:LYS:HD3	1:B:155:LYS:HA	1.70	0.46
2:C:906:PHE:HE2	3:D:1067:VAL:CA	2.19	0.46
1:A:55:SER:HB3	1:A:143:ARG:HB3	1.97	0.46
2:C:109:LYS:HE2	2:C:368:THR:HG22	1.97	0.46
5:F:285:GLU:HA	5:F:286:PRO:HD3	1.77	0.46
2:C:224:GLU:CD	2:C:224:GLU:H	2.24	0.46
3:D:818:ARG:HE	3:D:820:GLU:CD	2.24	0.46
3:D:1003:VAL:HG21	3:D:1041:LEU:HG	1.98	0.46
3:D:1264:GLU:OE2	3:D:1425:THR:OG1	2.33	0.46
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.69	0.46
3:D:1045[B]:MET:HE1	3:D:1057:VAL:HG23	1.98	0.46
1:B:5:LYS:HD3	1:B:7:LYS:HZ2	1.81	0.45
2:C:905:ILE:O	2:C:906:PHE:HB2	2.15	0.45
2:C:361:MET:SD	2:C:361:MET:N	2.90	0.45
2:C:693:GLU:HA	2:C:696:LYS:HD2	1.97	0.45
3:D:784:ASP:HB2	3:D:939:PHE:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1493:LYS:O	3:D:1497:GLU:HG2	2.16	0.45
3:D:207:PHE:HE2	5:F:98:GLU:HG2	1.81	0.45
1:A:11:PHE:O	1:B:228:PRO:HA	2.16	0.45
3:D:158:TYR:CE1	3:D:454:ALA:HB3	2.51	0.45
3:D:613:ARG:HG3	3:D:618:LEU:HD23	1.98	0.45
3:D:1057:VAL:HG12	3:D:1059:SER:H	1.82	0.45
3:D:1205:TYR:O	3:D:1366:LYS:HD3	2.16	0.45
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.98	0.45
1:A:99:LEU:HD23	1:A:114:PHE:CG	2.52	0.45
2:C:100:LEU:HD13	2:C:367:LEU:O	2.17	0.45
2:C:177:GLU:HG3	2:C:178:PRO:HD2	1.99	0.45
3:D:561:GLY:HA3	5:F:132:ARG:HD3	1.99	0.45
1:B:6:LEU:HD13	1:B:6:LEU:H	1.82	0.44
2:C:1038:TRP:CE2	3:D:1099:VAL:HG11	2.52	0.44
3:D:633:VAL:C	3:D:635:PRO:HD3	2.42	0.44
3:D:822:ALA:HB3	3:D:825:ALA:HB2	1.98	0.44
2:C:578:VAL:HG23	2:C:671:ASN:CG	2.42	0.44
2:C:617:ASP:HB2	2:C:619:ARG:HG2	1.98	0.44
3:D:12:LEU:HD21	3:D:104:PHE:CZ	2.52	0.44
3:D:415:VAL:HG13	3:D:419:ASP:HB2	1.99	0.44
5:F:144:ILE:HB	5:F:147:LEU:HD13	1.99	0.44
2:C:94:LEU:HD22	2:C:118:ILE:HD11	1.98	0.44
2:C:1110:ASP:OD2	2:C:1114:GLY:N	2.42	0.44
1:A:20:TYR:OH	1:A:198:ARG:HD2	2.18	0.44
2:C:63:GLY:HA2	2:C:102:HIS:HD2	1.82	0.44
2:C:76:PRO:HG3	2:C:120:LEU:HD12	2.00	0.44
2:C:1054:THR:OG1	2:C:1055:LEU:N	2.51	0.44
3:D:646:LYS:HB3	3:D:688:TRP:CZ3	2.52	0.44
3:D:1020:LEU:HB3	3:D:1035:ILE:HD12	2.00	0.44
2:C:35:PRO:HA	2:C:36:PRO:HD3	1.90	0.44
2:C:87:ASP:HA	2:C:131:GLY:HA3	1.99	0.44
2:C:693:GLU:HG3	2:C:697:ARG:HH12	1.81	0.44
2:C:784:ASP:OD1	2:C:784:ASP:N	2.50	0.44
3:D:1047:LYS:N	3:D:1051:GLU:O	2.34	0.44
6:G:15:DT:H2'	6:G:16:DC:C6	2.53	0.44
2:C:163:ILE:HG23	2:C:171:TRP:NE1	2.32	0.44
2:C:976:ASP:OD1	2:C:978:ARG:HD3	2.18	0.44
3:D:472:ALA:O	3:D:476:GLU:HG2	2.18	0.44
3:D:881:LEU:O	3:D:885:ILE:HG13	2.18	0.44
3:D:1286:THR:HB	3:D:1289:LYS:H	1.82	0.44
3:D:1305:LEU:HD13	3:D:1309:ALA:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1350:GLU:OE2	3:D:1357:ARG:NH1	2.40	0.44
2:C:205:GLU:O	2:C:209:ARG:HG2	2.17	0.44
3:D:230:TRP:CZ3	3:D:232:GLU:HG2	2.52	0.44
5:F:181:GLU:O	5:F:185:GLN:HG2	2.18	0.44
3:D:895:VAL:HG11	3:D:922:LEU:HD21	2.00	0.44
5:F:172:ARG:O	5:F:176:ILE:HG12	2.18	0.44
7:H:15:DT:C2	7:H:16:DC:C4	3.06	0.44
1:B:12:THR:HG22	1:B:13:VAL:N	2.32	0.43
2:C:769:PRO:HG3	3:D:65:ARG:HH12	1.83	0.43
5:F:166:LEU:HD13	5:F:170:HIS:HB3	1.99	0.43
1:A:215:VAL:HG13	1:B:222:LEU:HB3	2.00	0.43
2:C:1116:ALA:HB2	3:D:88:TYR:HB3	2.00	0.43
3:D:96:ALA:HB2	3:D:555:LYS:HG2	2.00	0.43
3:D:715:ALA:HB3	3:D:764:LEU:HA	2.00	0.43
3:D:1481:VAL:HA	4:E:18:ARG:HH21	1.82	0.43
3:D:63:TYR:OH	3:D:74:GLU:OE2	2.31	0.43
3:D:631:ILE:HD11	3:D:743:ASP:HB2	2.00	0.43
7:H:15:DT:C2	7:H:16:DC:C5	3.05	0.43
7:H:15:DT:O4	7:H:16:DC:N4	2.51	0.43
2:C:66:LEU:HD11	2:C:98:LEU:HB3	1.99	0.43
3:D:634:GLY:HA3	3:D:637:LEU:HD12	2.01	0.43
3:D:784:ASP:HB2	3:D:939:PHE:HE1	1.83	0.43
3:D:1018:ASN:HA	3:D:1019:PRO:HD3	1.88	0.43
4:E:83:ASP:OD1	4:E:83:ASP:N	2.51	0.43
6:G:4:DG:N2	6:G:5:DC:C2	2.86	0.43
2:C:580:MET:HB3	2:C:584:GLU:CD	2.43	0.43
2:C:745:ILE:HD12	2:C:801:VAL:O	2.18	0.43
2:C:926:PHE:HE1	2:C:929:ARG:HH21	1.66	0.43
2:C:1090:LYS:HD3	2:C:1090:LYS:HA	1.73	0.43
3:D:106:LYS:HA	3:D:106:LYS:HD2	1.86	0.43
5:F:153:PRO:HA	5:F:156:VAL:HG22	2.00	0.43
3:D:1383:ASP:HA	3:D:1384:PRO:HD3	1.81	0.43
1:B:176:ARG:HD3	3:D:884:ARG:NH2	2.34	0.43
2:C:418:LEU:CD2	7:H:14:DG:C8	2.91	0.43
3:D:760:ARG:O	3:D:764:LEU:HB2	2.19	0.43
5:F:188:ILE:HG12	5:F:224:VAL:HG21	2.00	0.43
2:C:86:LYS:HB2	2:C:88:LEU:HG	2.00	0.43
3:D:1350:GLU:O	3:D:1354:LYS:HG3	2.19	0.43
4:E:37:ASN:OD1	4:E:37:ASN:N	2.41	0.43
2:C:462:ASP:HB3	2:C:468:ARG:HD2	2.01	0.43
3:D:614:PHE:HA	3:D:618:LEU:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:761:ILE:HD12	4:E:20:THR:HA	2.01	0.43
3:D:1010:ASN:OD1	3:D:1014:ASN:ND2	2.51	0.43
4:E:59:ASN:O	4:E:63:TRP:HD1	2.01	0.43
1:A:101:LEU:HD21	1:A:109:VAL:HG11	2.01	0.42
3:D:229:ALA:HB1	3:D:245:LEU:HD12	2.00	0.42
3:D:989:TYR:OH	3:D:1052:THR:O	2.34	0.42
7:H:10:DA:H2''	7:H:11:DG:O5'	2.18	0.42
2:C:124:ASP:OD2	2:C:407:LYS:NZ	2.42	0.42
2:C:274:ARG:HG3	2:C:275:TYR:N	2.34	0.42
3:D:167:GLU:O	3:D:394:LEU:HD12	2.19	0.42
1:A:25:LEU:HB3	1:A:28:LEU:HD11	2.01	0.42
2:C:1009:SER:HB3	3:D:651:GLU:O	2.18	0.42
3:D:758:GLU:OE1	3:D:1476:THR:OG1	2.24	0.42
1:A:53:VAL:HG22	1:A:144:VAL:HG22	2.01	0.42
3:D:1038:LEU:O	3:D:1060:SER:HB2	2.19	0.42
6:G:4:DG:H1	7:H:24:DC:N4	2.11	0.42
2:C:807:ARG:H	2:C:807:ARG:HG2	1.70	0.42
3:D:90:MET:SD	3:D:521:PRO:HD3	2.58	0.42
3:D:762:GLN:HA	11:E:101:HOH:O	2.19	0.42
5:F:321:ILE:HA	5:F:321:ILE:HD13	1.78	0.42
2:C:878:SER:HA	3:D:1034:GLN:OE1	2.20	0.42
3:D:31:THR:HG21	5:F:257:THR:HG22	2.01	0.42
2:C:425:PHE:O	2:C:426:ASP:C	2.61	0.42
3:D:333:LEU:HB3	3:D:335:LEU:HD13	2.01	0.42
3:D:483:HIS:CE1	3:D:488:ARG:HD3	2.54	0.42
3:D:899:LEU:HD22	3:D:917:GLN:HB3	2.02	0.42
3:D:1065:LEU:HD23	3:D:1069:GLU:HB3	2.02	0.42
2:C:43:GLY:O	2:C:46:ALA:HB3	2.20	0.42
2:C:1066:ALA:O	2:C:1070:ILE:HG13	2.19	0.42
3:D:1147:ARG:HD3	3:D:1188:VAL:HG11	2.02	0.42
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.51	0.42
2:C:127:PHE:O	2:C:133:ASP:HA	2.20	0.42
2:C:281:LEU:HD13	2:C:305:PRO:HB2	2.02	0.42
3:D:128:TYR:OH	3:D:579:ASP:OD2	2.26	0.42
3:D:134:VAL:CG2	3:D:151:GLN:H	2.29	0.42
3:D:142:LEU:HB2	3:D:161:LEU:HD11	2.02	0.42
3:D:372:ASP:HA	3:D:373:PRO:HD3	1.96	0.42
3:D:1377:LYS:HE3	3:D:1378:TYR:CZ	2.55	0.42
2:C:704:HIS:CE1	2:C:1000:MET:HE1	2.55	0.41
2:C:721:ARG:HH22	2:C:785:VAL:HG11	1.85	0.41
2:C:1094:ALA:HA	3:D:518:PRO:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:729:HIS:HA	3:D:730:PRO:HD3	1.91	0.41
1:A:199:ILE:HB	1:A:207:PRO:HB3	2.02	0.41
2:C:884:GLN:HB2	2:C:992:MET:HE2	2.03	0.41
2:C:1005:MET:HE2	2:C:1005:MET:HB3	1.91	0.41
3:D:12:LEU:HD23	3:D:12:LEU:HA	1.84	0.41
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.53	0.41
2:C:1115:LEU:HB3	3:D:85:VAL:HG12	2.02	0.41
3:D:206:ARG:HA	3:D:206:ARG:HD2	1.91	0.41
5:F:94:LEU:HG	5:F:190:ALA:HB1	2.02	0.41
2:C:40:GLU:O	2:C:45:GLN:HG2	2.20	0.41
2:C:92:ALA:HB2	2:C:120:LEU:HD11	2.02	0.41
2:C:456:ALA:HB3	2:C:459:ALA:HB2	2.02	0.41
3:D:371:ILE:HG23	5:F:230:LYS:HD2	2.02	0.41
3:D:1444:THR:O	3:D:1448:THR:HG23	2.19	0.41
1:A:133:GLU:HG2	1:A:134:GLU:H	1.85	0.41
3:D:689:ASP:CG	4:E:51:LEU:HD11	2.46	0.41
3:D:106:LYS:O	3:D:586:ARG:NH1	2.53	0.41
3:D:964:LEU:HD23	3:D:964:LEU:HA	1.94	0.41
7:H:16:DC:C2	7:H:17:DA:C6	3.09	0.41
2:C:912:PRO:O	2:C:916:GLU:HG3	2.21	0.41
3:D:319:ALA:HA	3:D:337:LEU:HD23	2.02	0.41
3:D:1144:LEU:HD23	3:D:1144:LEU:HA	1.84	0.41
3:D:1302:GLU:OE1	3:D:1304:LYS:HE3	2.20	0.41
4:E:57:ASP:O	4:E:63:TRP:NE1	2.49	0.41
1:A:150:TYR:CE2	1:A:152:PRO:HG3	2.56	0.41
2:C:657:ASP:OD2	2:C:663:ASN:N	2.50	0.41
2:C:1019:GLN:HG2	2:C:1058:ASP:HB3	2.03	0.41
3:D:961:LYS:HE3	3:D:961:LYS:HB2	1.90	0.41
3:D:1450:ALA:HA	3:D:1455:LYS:HD2	2.02	0.41
5:F:120:THR:HG22	5:F:122:LEU:HD13	2.03	0.41
1:A:36:LEU:HD23	1:A:36:LEU:HA	1.95	0.41
1:B:90:LEU:HD21	1:B:121:GLU:HB2	2.02	0.41
2:C:141:HIS:HE1	2:C:144:PRO:HD3	1.85	0.41
2:C:172:ILE:HG21	2:C:184:MET:HE3	2.03	0.41
3:D:28:LYS:HA	3:D:29:PRO:HD3	1.88	0.41
3:D:222:GLY:HA2	3:D:333:LEU:O	2.21	0.41
3:D:1296:SER:N	3:D:1299:PHE:O	2.52	0.41
2:C:605:LYS:HB3	2:C:610:ARG:HH11	1.86	0.40
2:C:771:GLU:O	2:C:775:ARG:HB2	2.21	0.40
3:D:171:LEU:HD23	3:D:171:LEU:HA	1.86	0.40
3:D:841:TYR:HB2	3:D:864:VAL:HG13	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:234:LYS:HE2	5:F:234:LYS:HB3	1.96	0.40
5:F:156:VAL:O	5:F:160:ASP:HB2	2.21	0.40
1:B:124:ASN:OD1	1:B:124:ASN:N	2.53	0.40
2:C:428:ARG:CD	2:C:449:ILE:O	2.63	0.40
2:C:627:ARG:CZ	2:C:640:ARG:HG3	2.51	0.40
3:D:298:VAL:HG12	3:D:302:GLN:NE2	2.37	0.40
3:D:479:GLU:HA	3:D:482:LYS:HE2	2.04	0.40
3:D:911:LEU:O	3:D:915:VAL:HG23	2.21	0.40
3:D:1480:PHE:O	4:E:18:ARG:NH2	2.55	0.40
1:A:124:ASN:OD1	1:A:124:ASN:N	2.53	0.40
2:C:582:GLY:N	2:C:584:GLU:OE2	2.55	0.40
3:D:1264:GLU:CD	3:D:1425:THR:HG1	2.29	0.40
3:D:1439:SER:OG	3:D:1467:ILE:CD1	2.69	0.40
1:B:101:LEU:HD11	1:B:113:ASP:HB2	2.03	0.40
3:D:25:GLU:HB2	3:D:92:HIS:CE1	2.56	0.40
3:D:231:VAL:O	3:D:236:TYR:OH	2.39	0.40
3:D:1084:THR:O	3:D:1088:THR:HG23	2.22	0.40
3:D:1493:LYS:H	3:D:1493:LYS:HG2	1.63	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	229/315 (73%)	226 (99%)	3 (1%)	0	100	100
1	B	225/315 (71%)	218 (97%)	7 (3%)	0	100	100
2	C	1108/1119 (99%)	1077 (97%)	31 (3%)	0	100	100
3	D	1484/1524 (97%)	1447 (98%)	37 (2%)	0	100	100
4	E	92/99 (93%)	88 (96%)	4 (4%)	0	100	100
5	F	344/443 (78%)	339 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	3482/3815 (91%)	3395 (98%)	87 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	200/273 (73%)	196 (98%)	4 (2%)	48	68
1	B	198/273 (72%)	195 (98%)	3 (2%)	57	72
2	C	934/941 (99%)	902 (97%)	32 (3%)	32	59
3	D	1255/1279 (98%)	1226 (98%)	29 (2%)	44	66
4	E	82/88 (93%)	80 (98%)	2 (2%)	43	65
5	F	300/388 (77%)	294 (98%)	6 (2%)	48	68
All	All	2969/3242 (92%)	2893 (97%)	76 (3%)	40	64

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	66	SER
1	A	134	GLU
1	A	184	THR
1	B	5	LYS
1	B	6	LEU
1	B	154	GLU
2	C	64	LEU
2	C	177	GLU
2	C	182	VAL
2	C	200	LEU
2	C	261	ILE
2	C	274	ARG
2	C	284	ARG

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Mol	Chain	Res	Type
2	C	336	VAL
2	C	342	ASP
2	C	348	LEU
2	C	358	ARG
2	C	402	SER
2	C	419	THR
2	C	425	PHE
2	C	513	VAL
2	C	562	SER
2	C	583	LEU
2	C	584	GLU
2	C	610	ARG
2	C	640	ARG
2	C	657	ASP
2	C	670	GLN
2	C	715	THR
2	C	722	ILE
2	C	775	ARG
2	C	784	ASP
2	C	807	ARG
2	C	815	LEU
2	C	907	ASP
2	C	968	LEU
2	C	1014	SER
2	C	1060	ILE
3	D	35	ARG
3	D	40	GLU
3	D	183	GLU
3	D	231	VAL
3	D	335	LEU
3	D	354	VAL
3	D	427	VAL
3	D	452	ILE
3	D	471	GLU
3	D	530	VAL
3	D	618	LEU
3	D	632	VAL
3	D	687	VAL
3	D	754	PHE
3	D	810	GLU
3	D	894	LYS
3	D	1039	CYS

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Mol	Chain	Res	Type
3	D	1055	VAL
3	D	1067	VAL
3	D	1100	ASP
3	D	1129	THR
3	D	1188	VAL
3	D	1283	ILE
3	D	1307	LYS
3	D	1433	SER
3	D	1476	THR
3	D	1486	VAL
3	D	1487	VAL
3	D	1493	LYS
4	E	49	GLN
4	E	50	THR
5	F	122	LEU
5	F	141	VAL
5	F	154	LYS
5	F	205	ARG
5	F	321	ILE
5	F	338	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	ASN
1	B	95	GLN
2	C	117	HIS
2	C	219	GLN
2	C	498	GLN
2	C	647	GLN
2	C	1019	GLN
3	D	714	GLN
3	D	756	GLN
3	D	855	HIS
3	D	976	GLN
3	D	1116	ASN
3	D	1124	GLN
3	D	1334	GLN
5	F	170	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	3/4 (75%)	0	0

There are no RNA backbone outliers to report.

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 5 ligands modelled in this entry, 5 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%)	-0.20	0 100 100	63, 91, 117, 162	0
1	B	227/315 (72%)	-0.11	2 (0%) 81 65	63, 95, 124, 155	0
2	C	1112/1119 (99%)	-0.15	3 (0%) 90 82	45, 88, 139, 167	0
3	D	1486/1524 (97%)	-0.11	14 (0%) 81 65	46, 85, 145, 169	3 (0%)
4	E	94/99 (94%)	-0.07	0 100 100	65, 97, 135, 142	0
5	F	346/443 (78%)	0.04	3 (0%) 81 65	65, 103, 147, 164	0
6	G	17/21 (80%)	-0.16	0 100 100	70, 105, 177, 184	0
7	H	24/27 (88%)	0.12	0 100 100	96, 126, 176, 188	0
8	I	4/4 (100%)	-0.86	0 100 100	69, 72, 79, 89	0
All	All	3541/3867 (91%)	-0.11	22 (0%) 85 73	45, 91, 142, 188	3 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	184	GLU	3.4
3	D	203	ALA	3.2
3	D	353	VAL	3.0
3	D	560	GLN	2.8
3	D	292	VAL	2.8
3	D	420	VAL	2.7
3	D	290	PRO	2.7
2	C	64	LEU	2.7
3	D	368	VAL	2.7
3	D	303	PRO	2.6
1	B	2	LEU	2.6
2	C	188	LYS	2.5
3	D	291	LEU	2.5
5	F	141	VAL	2.4
3	D	1252	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
5	F	146	GLY	2.2
5	F	145	PRO	2.2
3	D	409	VAL	2.1
2	C	63	GLY	2.1
3	D	452	ILE	2.1
1	B	91	ASN	2.0
3	D	297	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	D	2004	1/1	0.85	0.23	63,63,63,63	0
9	MG	B	2001	1/1	0.95	0.05	95,95,95,95	0
9	MG	D	2003	1/1	0.98	0.05	52,52,52,52	0
10	ZN	D	2001	1/1	0.99	0.04	71,71,71,71	0
10	ZN	D	2002	1/1	0.99	0.04	111,111,111,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.