



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:30 AM UTC

PDB ID : 6KQM / pdb_00006kqm
Title : Thermus thermophilus initial transcription complex comprising sigma A and 5'-triphosphate RNA of 5 nt
Authors : Zhang, Y.; Li, L.; Ebright, R.H.
Deposited on : 2019-08-18
Resolution : 3.20 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
Mogul : 2022.3.0, CSD as543be (2022)
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

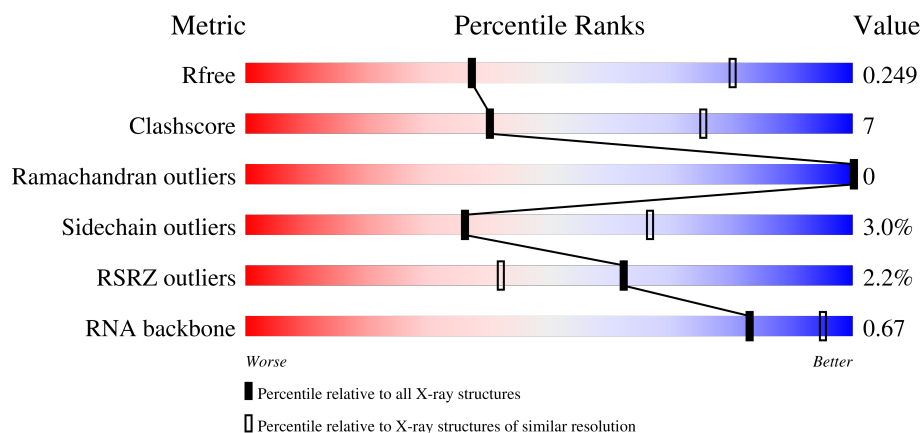
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)
RNA backbone	3983	1222 (3.50-2.90)

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	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	E	99	% 77% 17% • 5%
5	F	443	8% 65% 12% • 22%
6	G	21	5% 57% 24% 19%
7	H	27	4% 56% 33% 11%
8	I	5	80% 20%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 28789 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	3	0
			11753	7455	2067	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
			761	486	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
			2807	1770	509	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(*(CTP))-R(P*UP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	I	5	Total	C	N	O	P	0	0	0
			106	47	18	36	5			

- 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

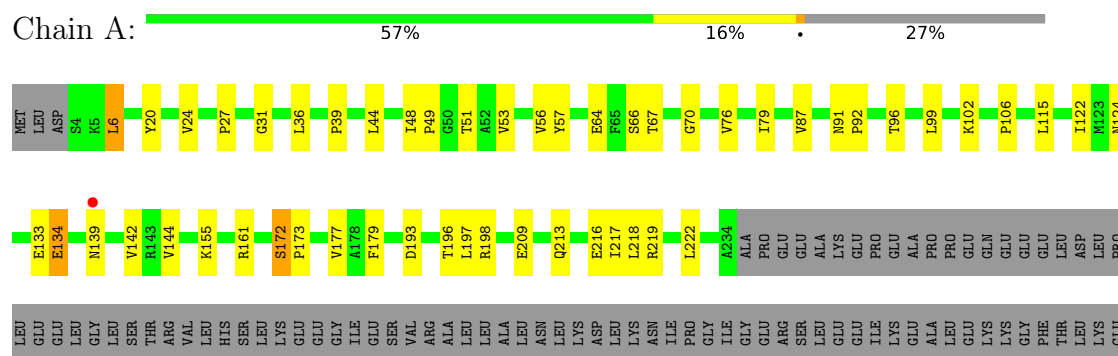
- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	12	Total O 12 12	0	0
11	B	6	Total O 6 6	0	0
11	C	44	Total O 44 44	0	0
11	D	61	Total O 61 61	0	0
11	E	2	Total O 2 2	0	0
11	F	6	Total O 6 6	0	0
11	G	9	Total O 9 9	0	0
11	H	2	Total O 2 2	0	0
11	I	5	Total O 5 5	0	0

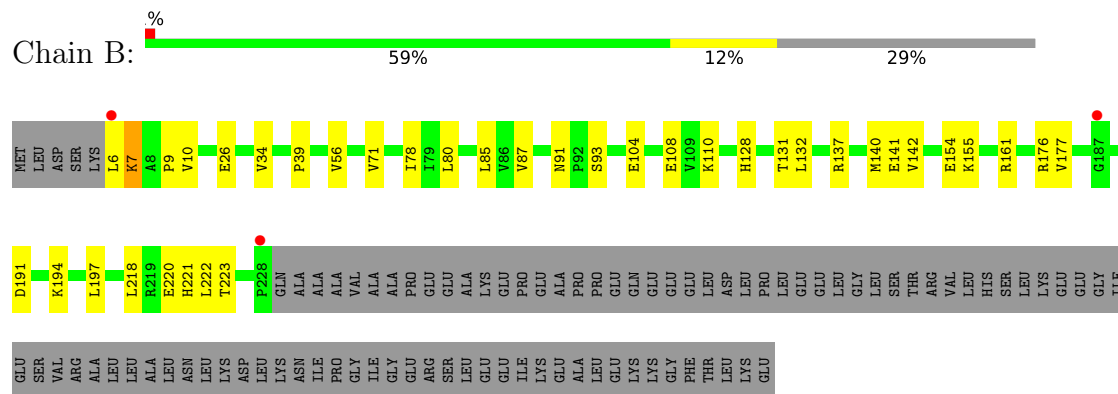
3 Residue-property plots [i](#)

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.

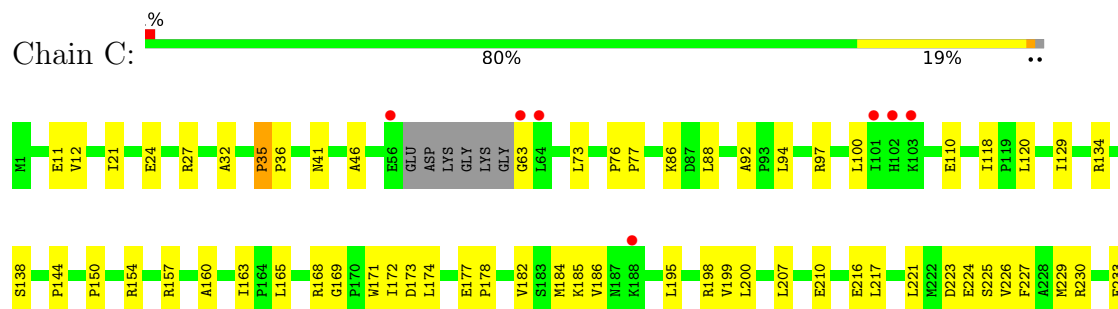
- Molecule 1: DNA-directed RNA polymerase subunit alpha

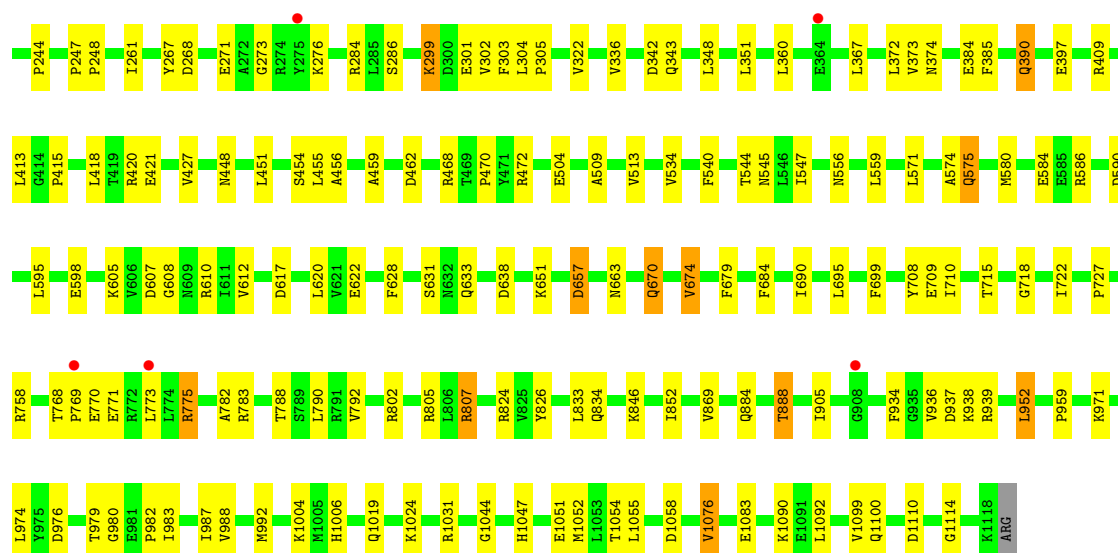


- Molecule 1: DNA-directed RNA polymerase subunit alpha

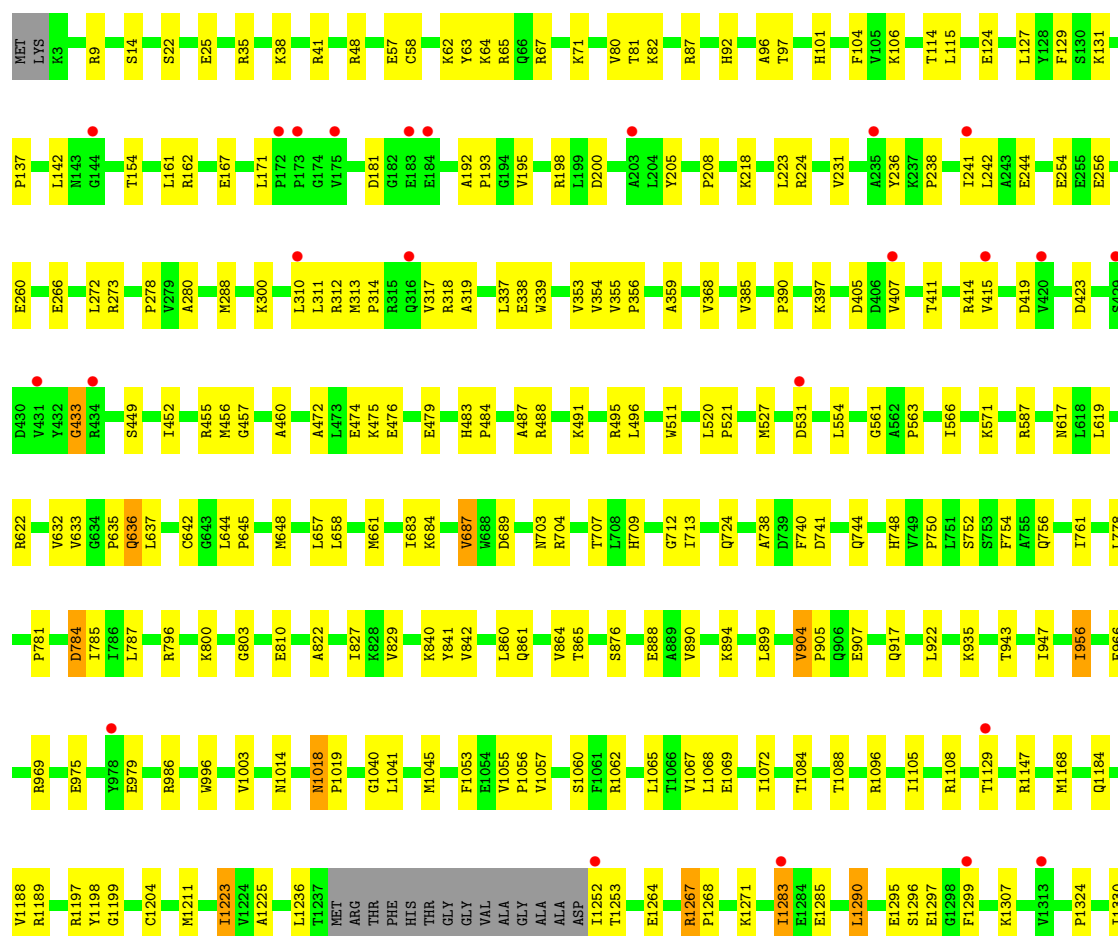
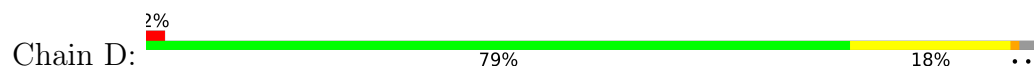


- Molecule 2: DNA-directed RNA polymerase subunit beta

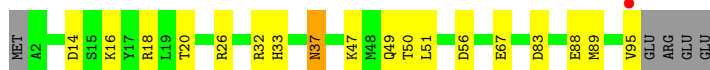
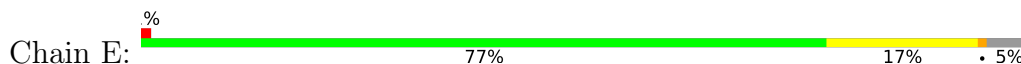




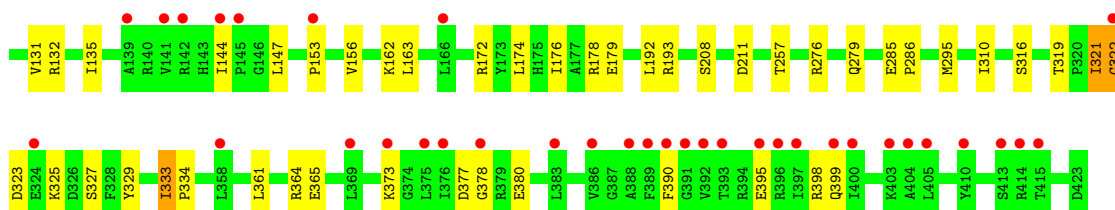
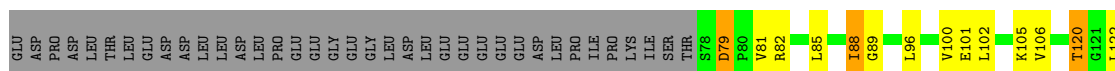
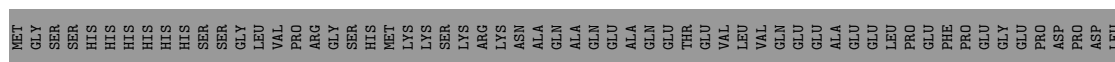
• Molecule 3: DNA-directed RNA polymerase subunit beta'



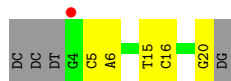
- Molecule 4: DNA-directed RNA polymerase subunit omega



- Molecule 5: RNA polymerase sigma factor SigA



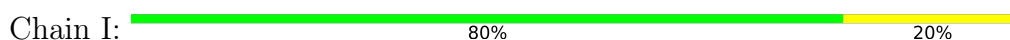
- Molecule 6: DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*CP*GP*AP*GP*GP*G)-3')



- 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(* (CTP))-R(P*UP*CP*GP*A)-3')





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	184.73Å 102.75Å 295.71Å 90.00° 98.86° 90.00°	Depositor
Resolution (Å)	48.89 – 3.20 48.89 – 3.20	Depositor EDS
% Data completeness (in resolution range)	95.1 (48.89-3.20) 95.0 (48.89-3.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.205 , 0.252 0.204 , 0.249	Depositor DCC
R_{free} test set	4310 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	28789	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: MG, ZN, CTP

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	0.82	4/2515 (0.2%)
1	B	0.33	0/1790	0.83	1/2435 (0.0%)
2	C	0.32	0/8969	0.79	9/12129 (0.1%)
3	D	0.35	2/11969 (0.0%)	0.79	10/16182 (0.1%)
4	E	0.30	0/775	0.76	0/1045
5	F	0.31	0/2852	0.77	5/3837 (0.1%)
6	G	0.19	0/393	0.53	0/606
7	H	0.21	0/556	0.54	0/858
8	I	0.14	0/94	0.26	0/144
All	All	0.33	2/29247 (0.0%)	0.78	29/39751 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1223	ILE	C-O	-8.25	1.14	1.24
3	D	1225	ALA	C-O	-7.69	1.15	1.24

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	433	GLY	N-CA-C	7.50	120.67	110.69
2	C	169	GLY	CA-C-N	7.03	126.95	119.85
2	C	169	GLY	C-N-CA	7.03	126.95	119.85
1	A	172	SER	CA-C-N	6.33	125.95	119.56
1	A	172	SER	C-N-CA	6.33	125.95	119.56
3	D	1014	ASN	N-CA-C	6.09	119.18	111.69
5	F	322	GLY	N-CA-C	-6.04	103.92	112.37
2	C	271	GLU	N-CA-C	-6.00	104.74	111.28
5	F	323	ASP	N-CA-C	5.99	119.04	110.24
3	D	1340	GLY	CA-C-N	5.96	125.17	118.97

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1340	GLY	C-N-CA	5.96	125.17	118.97
2	C	35	PRO	CA-C-N	5.73	126.12	119.47
2	C	35	PRO	C-N-CA	5.73	126.12	119.47
3	D	740	PHE	N-CA-C	5.59	119.73	113.02
3	D	617	ASN	N-CA-C	5.57	120.19	113.17
2	C	792	VAL	CA-C-N	5.56	123.72	119.66
2	C	792	VAL	C-N-CA	5.56	123.72	119.66
3	D	1481	VAL	N-CA-C	5.36	118.14	112.83
3	D	1018	ASN	CA-C-N	5.31	124.76	119.24
3	D	1018	ASN	C-N-CA	5.31	124.76	119.24
3	D	740	PHE	CB-CA-C	5.29	119.78	110.37
5	F	333	ILE	N-CA-C	5.28	112.78	107.55
1	B	9	PRO	CB-CA-C	-5.25	103.35	110.60
2	C	980	GLY	N-CA-C	-5.24	107.30	114.64
5	F	79	ASP	CA-C-N	5.21	124.83	119.05
5	F	79	ASP	C-N-CA	5.21	124.83	119.05
1	A	115	LEU	CA-C-N	5.15	125.14	119.89
1	A	115	LEU	C-N-CA	5.15	125.14	119.89
2	C	303	PHE	N-CA-C	5.07	117.19	111.11

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	34	0
1	B	1758	0	1808	26	0
2	C	8792	0	8902	130	0
3	D	11753	0	11992	157	0
4	E	761	0	778	16	0
5	F	2807	0	2882	42	0
6	G	350	0	192	4	0
7	H	495	0	272	8	0
8	I	106	0	54	1	0
9	B	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	D	2	0	0	0	0
9	F	1	0	0	0	0
10	D	2	0	0	0	0
11	A	12	0	0	0	0
11	B	6	0	0	0	0
11	C	44	0	0	1	0
11	D	61	0	0	0	0
11	E	2	0	0	0	0
11	F	6	0	0	1	0
11	G	9	0	0	0	0
11	H	2	0	0	0	0
11	I	5	0	0	0	0
All	All	28789	0	28749	376	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 (376) 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)
4:E:33:HIS:HA	4:E:95:VAL:CG2	1.78	1.13
4:E:33:HIS:HA	4:E:95:VAL:HG21	1.48	0.92
2:C:768:THR:CG2	2:C:769:PRO:HD2	2.06	0.85
2:C:628:PHE:H	2:C:638:ASP:HB3	1.41	0.85
4:E:33:HIS:HA	4:E:95:VAL:HG22	1.54	0.85
2:C:768:THR:HG22	2:C:770:GLU:OE1	1.82	0.80
2:C:12:VAL:HG21	2:C:472:ARG:HD3	1.63	0.80
4:E:32:ARG:O	4:E:95:VAL:HG11	1.85	0.77
5:F:82:ARG:HB2	7:H:8:DG:O6	1.86	0.76
2:C:768:THR:HG23	2:C:769:PRO:HD2	1.67	0.76
2:C:168:ARG:HD3	2:C:268:ASP:HB3	1.68	0.75
7:H:18:DC:H2''	7:H:19:DG:C8	2.23	0.74
2:C:165:LEU:HB2	2:C:168:ARG:HG3	1.72	0.71
2:C:768:THR:HG22	2:C:769:PRO:HD2	1.73	0.70
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.26	0.68
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.76	0.68
4:E:33:HIS:HE2	4:E:89:MET:HB3	1.59	0.67
7:H:12:DC:O5'	7:H:12:DC:H6	1.78	0.67
3:D:65:ARG:NH1	5:F:378:GLY:O	2.29	0.66
3:D:256:GLU:HG3	3:D:300:LYS:HG3	1.78	0.66
1:A:67:THR:HG21	2:C:608:GLY:HA3	1.76	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:566:ILE:HD11	5:F:192:LEU:HD21	1.79	0.65
2:C:24:GLU:OE2	2:C:27:ARG:NH2	2.30	0.65
1:B:6:LEU:HD12	1:B:6:LEU:N	2.12	0.65
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.78	0.65
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.80	0.64
3:D:784:ASP:N	3:D:784:ASP:OD1	2.31	0.64
3:D:208:PRO:HA	3:D:390:PRO:HA	1.79	0.64
5:F:321:ILE:HD11	5:F:327:SER:HB3	1.78	0.64
2:C:805:ARG:O	2:C:807[A]:ARG:NH2	2.28	0.63
2:C:63:GLY:HA3	2:C:100:LEU:HD21	1.79	0.63
2:C:758:ARG:HH21	2:C:788:THR:HB	1.62	0.63
3:D:899:LEU:HD22	3:D:917:GLN:HB3	1.81	0.63
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.78	0.63
5:F:316:SER:O	5:F:319:THR:HG23	1.98	0.63
3:D:273:ARG:HB3	3:D:278:PRO:HA	1.81	0.63
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.80	0.62
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.33	0.62
3:D:657:LEU:HG	3:D:661:MET:HE2	1.81	0.62
2:C:1019:GLN:HG2	2:C:1058:ASP:HB3	1.81	0.62
1:A:56:VAL:HG22	1:A:142:VAL:HG12	1.82	0.62
3:D:1236:LEU:HA	3:D:1359:GLN:HG3	1.82	0.62
1:B:7:LYS:HD3	1:B:7:LYS:N	2.15	0.61
2:C:1083:GLU:OE2	3:D:87:ARG:NH2	2.34	0.61
2:C:768:THR:CG2	2:C:769:PRO:CD	2.79	0.60
2:C:420:ARG:NH1	8:I:1:CTP:O2A	2.34	0.60
3:D:411:THR:O	5:F:178:ARG:NH1	2.33	0.60
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.83	0.60
3:D:405:ASP:HB3	3:D:423:ASP:HA	1.84	0.59
2:C:129:ILE:HB	2:C:134:ARG:HD2	1.84	0.59
2:C:657:ASP:OD2	2:C:663:ASN:N	2.35	0.59
3:D:272:LEU:HB2	3:D:280:ALA:HB3	1.83	0.59
2:C:226:VAL:HA	2:C:229:MET:HE2	1.84	0.59
5:F:319:THR:OG1	5:F:329:TYR:HB2	2.04	0.58
3:D:787:LEU:HD21	3:D:947:ILE:HG21	1.85	0.58
2:C:462:ASP:HB3	2:C:468:ARG:HD2	1.86	0.57
3:D:319:ALA:HA	3:D:337:LEU:HD23	1.87	0.57
2:C:172:ILE:HG12	2:C:186:VAL:HG22	1.87	0.57
3:D:644:LEU:HD12	3:D:645:PRO:HD2	1.85	0.57
1:A:31:GLY:N	1:A:193:ASP:OD1	2.38	0.57
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.87	0.56
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:802:ARG:HB2	2:C:826:TYR:HB2	1.87	0.56
3:D:890:VAL:HB	3:D:922:LEU:HD13	1.87	0.56
3:D:1096:ARG:NH1	3:D:1440:PHE:O	2.38	0.56
2:C:884:GLN:O	2:C:888:THR:OG1	2.24	0.56
3:D:236:TYR:HB2	3:D:319:ALA:HB3	1.87	0.56
1:A:106:PRO:HG3	1:A:134:GLU:HG2	1.86	0.56
1:B:80:LEU:HD21	3:D:842:VAL:HG12	1.88	0.55
3:D:241:ILE:HA	3:D:312:ARG:HG2	1.88	0.55
2:C:41:ASN:O	2:C:46:ALA:HB2	2.06	0.55
2:C:768:THR:CG2	2:C:770:GLU:OE1	2.53	0.55
3:D:167:GLU:OE2	3:D:198:ARG:NH1	2.40	0.55
1:A:133:GLU:OE1	2:C:610:ARG:NH1	2.40	0.55
2:C:575:GLN:HG2	2:C:670:GLN:HE21	1.72	0.55
2:C:905:ILE:HG13	2:C:905:ILE:O	2.07	0.55
2:C:160:ALA:HB3	2:C:174:LEU:HB2	1.89	0.54
2:C:775:ARG:HD3	2:C:782:ALA:HB2	1.88	0.54
3:D:637:LEU:HD13	3:D:642:CYS:HA	1.90	0.54
3:D:904:VAL:HG22	3:D:905:PRO:HD2	1.88	0.54
3:D:142:LEU:HB2	3:D:161:LEU:HD21	1.89	0.54
4:E:32:ARG:O	4:E:95:VAL:CG1	2.54	0.54
1:A:36:LEU:HD11	1:B:221:HIS:HB3	1.88	0.54
4:E:14:ASP:OD2	4:E:18:ARG:NH1	2.40	0.54
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.40	0.54
2:C:207:LEU:HD13	2:C:221:LEU:HD21	1.90	0.54
3:D:1462:LEU:HD22	3:D:1472:ILE:HB	1.91	0.53
5:F:79:ASP:OD2	7:H:8:DG:N1	2.30	0.53
3:D:171:LEU:HD12	3:D:390:PRO:HG2	1.90	0.53
2:C:97:ARG:NH1	2:C:110:GLU:OE1	2.35	0.53
5:F:321:ILE:HG12	5:F:322:GLY:H	1.74	0.53
2:C:612:VAL:HG22	2:C:622:GLU:HG3	1.91	0.53
3:D:1045[B]:MET:HE1	3:D:1057:VAL:HG23	1.91	0.53
5:F:279:GLN:HB3	5:F:286:PRO:HD3	1.91	0.53
3:D:761:ILE:HD12	4:E:20:THR:HA	1.91	0.52
3:D:318:ARG:NH1	3:D:338:GLU:OE1	2.42	0.52
1:B:78:ILE:HG21	1:B:140:MET:HE1	1.91	0.52
2:C:715:THR:OG1	2:C:718:GLY:O	2.25	0.52
2:C:390:GLN:HB3	2:C:415:PRO:HD3	1.90	0.52
2:C:468:ARG:NH2	11:C:1205:HOH:O	2.43	0.52
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.40	0.52
3:D:223:LEU:HD11	3:D:288:MET:HE1	1.92	0.52
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:367:LEU:HD13	2:C:372:LEU:HD21	1.92	0.52
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.10	0.52
1:B:6:LEU:N	1:B:6:LEU:CD1	2.73	0.51
2:C:605:LYS:HB2	2:C:612:VAL:HB	1.92	0.51
5:F:193:ARG:HB2	7:H:6:DT:H1'	1.92	0.51
1:B:7:LYS:HD3	1:B:7:LYS:H	1.75	0.51
1:B:26:GLU:HB3	1:B:194:LYS:HG3	1.93	0.51
2:C:574:ALA:HA	2:C:670:GLN:HE22	1.74	0.51
3:D:561:GLY:HA3	5:F:132:ARG:HD3	1.92	0.51
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.91	0.51
3:D:781:PRO:HB3	3:D:785:ILE:CG2	2.41	0.51
2:C:1110:ASP:OD2	2:C:1114:GLY:N	2.31	0.51
3:D:1296:SER:OG	3:D:1297:GLU:N	2.43	0.51
2:C:1004:LYS:HD3	3:D:744:GLN:NE2	2.26	0.51
2:C:690:ILE:HG13	2:C:852:ILE:HG23	1.93	0.51
3:D:956:ILE:HD11	3:D:1062:ARG:HG2	1.91	0.51
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.10	0.51
1:B:220:GLU:O	1:B:223:THR:OG1	2.27	0.50
2:C:177:GLU:HG3	2:C:178:PRO:HD2	1.91	0.50
3:D:244:GLU:HG3	3:D:310:LEU:HG	1.92	0.50
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.92	0.50
3:D:142:LEU:HD13	3:D:161:LEU:HD11	1.93	0.50
2:C:773:LEU:HD13	5:F:373:LYS:HG3	1.93	0.50
1:A:53:VAL:HG22	1:A:144:VAL:HG22	1.93	0.50
3:D:96:ALA:HB3	3:D:554:LEU:HD23	1.93	0.50
3:D:703:ASN:HB2	3:D:713:ILE:HG12	1.94	0.50
3:D:181:ASP:HB2	3:D:205:TYR:CD1	2.46	0.50
4:E:32:ARG:O	4:E:95:VAL:HG21	2.12	0.50
6:G:5:DC:H2''	6:G:6:DA:OP2	2.12	0.50
2:C:420:ARG:NH2	2:C:448:ASN:OD1	2.44	0.49
2:C:1006:HIS:HB2	2:C:1024:LYS:HG3	1.94	0.49
3:D:63:TYR:HB2	3:D:80:VAL:HG21	1.94	0.49
3:D:224:ARG:NE	3:D:254:GLU:OE2	2.37	0.49
1:A:222:LEU:HD11	1:B:218:LEU:HG	1.95	0.49
1:A:218:LEU:HD23	1:B:222:LEU:HD21	1.94	0.49
2:C:768:THR:HG22	2:C:769:PRO:CD	2.41	0.49
3:D:1168:MET:HE3	3:D:1168:MET:HA	1.95	0.49
1:B:104:GLU:OE2	1:B:137:ARG:NH1	2.46	0.49
2:C:224:GLU:CD	2:C:224:GLU:H	2.20	0.49
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.93	0.49
3:D:996:TRP:CD2	3:D:1056:PRO:HG3	2.48	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:153:PRO:HA	5:F:156:VAL:HG22	1.95	0.49
2:C:937:ASP:OD1	2:C:938:LYS:N	2.45	0.49
2:C:168:ARG:O	2:C:267:TYR:HA	2.13	0.48
2:C:679:PHE:HA	3:D:943:THR:HG23	1.95	0.48
3:D:14:SER:HB3	3:D:511:TRP:CE2	2.48	0.48
3:D:67:ARG:HB3	5:F:377:ASP:O	2.13	0.48
1:A:216:GLU:OE2	1:A:219:ARG:NH2	2.47	0.48
3:D:619:LEU:HD11	3:D:1439:SER:HB2	1.94	0.48
1:B:191:ASP:OD1	1:B:191:ASP:N	2.44	0.48
5:F:321:ILE:HD13	5:F:321:ILE:H	1.78	0.48
7:H:2:DA:C8	7:H:2:DA:H5'	2.48	0.48
2:C:92:ALA:HB2	2:C:120:LEU:HD11	1.95	0.48
2:C:727:PRO:HB3	2:C:783:ARG:HD3	1.94	0.48
3:D:127:LEU:HA	3:D:457:GLY:HA2	1.96	0.48
3:D:353:VAL:HG22	3:D:368:VAL:HG22	1.96	0.48
1:A:6:LEU:HD11	1:A:27:PRO:HG2	1.96	0.48
3:D:1264:GLU:OE2	3:D:1425:THR:OG1	2.31	0.48
2:C:351:LEU:HD11	2:C:373:VAL:HG13	1.96	0.48
2:C:547:ILE:O	2:C:905:ILE:CG2	2.62	0.48
5:F:319:THR:OG1	5:F:329:TYR:CB	2.62	0.48
3:D:22:SER:HB2	3:D:92:HIS:HB3	1.95	0.48
1:A:39:PRO:HG3	1:B:39:PRO:HG3	1.96	0.47
2:C:244:PRO:O	5:F:82:ARG:NH1	2.47	0.47
2:C:710:ILE:HD12	2:C:790:LEU:HB2	1.95	0.47
5:F:120:THR:HG22	5:F:122:LEU:HD13	1.94	0.47
1:A:133:GLU:HG2	1:A:134:GLU:H	1.79	0.47
3:D:455:ARG:HB2	3:D:460:ALA:HB2	1.96	0.47
3:D:1057:VAL:HG22	3:D:1069:GLU:HG2	1.95	0.47
3:D:1105:ILE:HG23	3:D:1199:GLY:HA2	1.96	0.47
5:F:172:ARG:O	5:F:176:ILE:HG12	2.14	0.47
3:D:314:PRO:HB2	3:D:317:VAL:HG12	1.96	0.47
1:A:209:GLU:O	1:A:213:GLN:HG2	2.14	0.47
2:C:769:PRO:HG3	3:D:65:ARG:HH12	1.78	0.47
1:B:176:ARG:NH2	3:D:888:GLU:OE1	2.42	0.47
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.97	0.47
2:C:540:PHE:HB3	2:C:544:THR:HB	1.97	0.47
3:D:97:THR:HG21	3:D:571:LYS:HG2	1.96	0.47
2:C:580:MET:HB3	2:C:584:GLU:CD	2.40	0.47
1:B:110:LYS:HD3	1:B:128:HIS:HA	1.97	0.47
2:C:173:ASP:HB2	2:C:185:LYS:HB3	1.97	0.46
5:F:163:LEU:HD13	5:F:174:LEU:HD13	1.96	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:26:ARG:NE	4:E:67:GLU:OE1	2.44	0.46
2:C:21:ILE:HD12	2:C:455:LEU:HD22	1.98	0.46
2:C:547:ILE:O	2:C:905:ILE:HG22	2.15	0.46
2:C:144:PRO:HB2	2:C:273:GLY:HA3	1.96	0.46
3:D:1147:ARG:NH2	3:D:1369:GLU:OE1	2.45	0.46
4:E:47:LYS:NZ	4:E:56:ASP:OD1	2.39	0.46
5:F:88:ILE:HG23	5:F:193:ARG:HG2	1.97	0.46
1:A:64:GLU:HG3	1:A:79:ILE:HD12	1.97	0.46
2:C:456:ALA:HB3	2:C:459:ALA:HB2	1.98	0.46
2:C:598:GLU:O	2:C:651:LYS:NZ	2.45	0.46
3:D:433:GLY:HA2	3:D:449:SER:H	1.81	0.46
2:C:1052:MET:HE1	3:D:748:HIS:HB3	1.97	0.46
2:C:1090:LYS:HD3	2:C:1090:LYS:HA	1.73	0.46
3:D:483:HIS:CE1	3:D:488:ARG:HD3	2.51	0.46
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.69	0.46
2:C:695:LEU:HD21	2:C:833:LEU:HB3	1.97	0.46
5:F:333:ILE:HA	5:F:334:PRO:HD3	1.82	0.46
2:C:27:ARG:HB3	2:C:27:ARG:HH21	1.81	0.46
3:D:415:VAL:HG13	3:D:419:ASP:HB2	1.97	0.46
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.98	0.45
1:B:177:VAL:HG13	1:B:197:LEU:HD11	1.98	0.45
1:A:44:LEU:HB3	1:A:177:VAL:HG21	1.98	0.45
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.97	0.45
3:D:218:LYS:HG2	3:D:338:GLU:HG2	1.99	0.45
3:D:864:VAL:HG22	3:D:865:THR:H	1.81	0.45
5:F:361:LEU:HB3	5:F:365:GLU:HG3	1.98	0.45
3:D:114:THR:HG23	3:D:495:ARG:HG2	1.98	0.45
3:D:527:MET:HE2	3:D:527:MET:HB2	1.86	0.45
4:E:83:ASP:OD1	4:E:83:ASP:N	2.49	0.45
1:B:155:LYS:HA	1:B:155:LYS:HD3	1.69	0.45
2:C:32:ALA:HB2	2:C:73:LEU:HD12	1.98	0.45
5:F:316:SER:HB3	5:F:319:THR:HG23	1.97	0.45
3:D:129:PHE:CD1	3:D:456:MET:HB3	2.52	0.45
3:D:162:ARG:O	3:D:449:SER:HB2	2.17	0.45
3:D:840:LYS:HE3	3:D:841:TYR:CZ	2.52	0.45
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.99	0.45
2:C:374:ASN:OD1	5:F:276:ARG:HD2	2.17	0.45
2:C:1054:THR:OG1	2:C:1055:LEU:N	2.49	0.45
2:C:556:ASN:O	2:C:559:LEU:HB3	2.17	0.44
3:D:966:GLU:O	3:D:969:ARG:HG2	2.17	0.44
1:A:91:ASN:HA	1:A:92:PRO:HD3	1.83	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:470:PRO:HB2	2:C:534:VAL:HG21	1.99	0.44
2:C:674:VAL:HG21	2:C:992:MET:HE3	1.98	0.44
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.98	0.44
5:F:102:LEU:O	5:F:106:VAL:HG23	2.18	0.44
2:C:952:LEU:HD12	2:C:952:LEU:HA	1.78	0.44
3:D:162:ARG:O	3:D:414:ARG:NH1	2.51	0.44
2:C:939:ARG:HG2	2:C:982:PRO:HD3	2.00	0.44
2:C:1092:LEU:HD13	2:C:1099:VAL:HG21	2.00	0.44
3:D:1003:VAL:HG21	3:D:1041:LEU:HG	2.00	0.44
2:C:768:THR:HG23	2:C:769:PRO:CD	2.44	0.44
3:D:860:LEU:O	3:D:876:SER:HB2	2.18	0.44
3:D:1324:PRO:HG3	3:D:1330:ILE:HD11	1.98	0.44
5:F:135:ILE:HD11	5:F:178:ARG:HB3	2.00	0.44
1:A:124:ASN:OD1	1:A:124:ASN:N	2.50	0.44
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.52	0.44
3:D:101:HIS:HB3	3:D:104:PHE:HD2	1.82	0.44
1:A:51:THR:OG1	1:A:87:VAL:O	2.31	0.44
2:C:172:ILE:HD13	2:C:184:MET:HE3	1.99	0.44
5:F:322:GLY:HA3	6:G:20:DG:N2	2.33	0.44
1:A:57:TYR:CG	1:A:161:ARG:HD2	2.53	0.43
1:A:99:LEU:HD21	1:A:122:ILE:HD11	1.99	0.43
3:D:487:ALA:O	3:D:491:LYS:HG2	2.18	0.43
2:C:154:ARG:HH21	2:C:157:ARG:HG3	1.83	0.43
3:D:935:LYS:HE2	3:D:935:LYS:HB3	1.90	0.43
3:D:1065:LEU:HD23	3:D:1069:GLU:HB3	2.00	0.43
3:D:1353:GLN:O	3:D:1357:ARG:HG3	2.18	0.43
3:D:1364:HIS:CE1	3:D:1366:LYS:HE2	2.53	0.43
5:F:89:GLY:HA3	7:H:7:DG:C6	2.54	0.43
2:C:150:PRO:HG3	2:C:322:VAL:HG11	1.99	0.43
3:D:1375:MET:HE2	3:D:1421:LEU:HD11	2.01	0.43
5:F:101:GLU:HG2	5:F:105:LYS:HE2	2.00	0.43
3:D:633:VAL:C	3:D:635:PRO:HD3	2.43	0.43
5:F:257:THR:N	11:F:2101:HOH:O	2.50	0.43
1:A:70:GLY:N	2:C:607:ASP:OD1	2.51	0.43
1:A:20:TYR:OH	1:A:198:ARG:HD2	2.18	0.43
3:D:41:ARG:HE	3:D:48:ARG:CZ	2.32	0.43
3:D:131:LYS:NZ	3:D:154:THR:HG22	2.34	0.43
6:G:5:DC:C2	6:G:6:DA:N7	2.87	0.43
2:C:76:PRO:HG3	2:C:120:LEU:HD12	2.00	0.43
3:D:161:LEU:HB3	3:D:452:ILE:HD11	2.01	0.43
5:F:395:GLU:OE2	5:F:398:ARG:NH2	2.52	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:708:TYR:HB3	2:C:790:LEU:HD21	2.01	0.43
3:D:707:THR:HG23	3:D:712:GLY:HA3	2.00	0.43
3:D:520:LEU:HD12	3:D:521:PRO:HD2	2.00	0.43
3:D:1018:ASN:HA	3:D:1019:PRO:HD3	1.88	0.43
3:D:1211:MET:HE1	4:E:16:LYS:HD2	2.01	0.43
3:D:1285:GLU:HG3	3:D:1290:LEU:HG	2.00	0.43
3:D:1383:ASP:HA	3:D:1384:PRO:HD3	1.79	0.43
3:D:1455:LYS:HE3	3:D:1455:LYS:HB2	1.88	0.43
3:D:57:GLU:HG3	3:D:64:LYS:HG2	2.01	0.42
3:D:483:HIS:CG	3:D:484:PRO:HD2	2.55	0.42
5:F:81:VAL:O	5:F:85:LEU:HG	2.20	0.42
2:C:35:PRO:HA	2:C:36:PRO:HD3	1.87	0.42
2:C:1047:HIS:O	2:C:1051:GLU:HB2	2.19	0.42
3:D:683:ILE:HG23	3:D:687:VAL:HG21	2.01	0.42
3:D:800:LYS:HB3	3:D:822:ALA:HB2	2.01	0.42
3:D:975:GLU:O	3:D:979:GLU:HG2	2.19	0.42
2:C:684:PHE:HB3	3:D:633:VAL:HG21	2.01	0.42
3:D:704:ARG:HD2	3:D:738:ALA:HB2	2.01	0.42
2:C:846:LYS:HE3	3:D:741:ASP:HB2	2.01	0.42
3:D:25:GLU:HB2	3:D:92:HIS:CE1	2.55	0.42
3:D:840:LYS:HE3	3:D:841:TYR:OH	2.19	0.42
3:D:1336:LEU:HB2	3:D:1344:VAL:HG21	2.01	0.42
5:F:88:ILE:HD11	5:F:192:LEU:HD13	2.00	0.42
3:D:353:VAL:HG12	3:D:355:VAL:H	1.84	0.42
3:D:563:PRO:HD2	3:D:566:ILE:HD12	2.00	0.42
3:D:689:ASP:OD2	4:E:51:LEU:HD11	2.20	0.42
1:B:7:LYS:N	1:B:7:LYS:CD	2.73	0.42
7:H:3:DT:H2'	7:H:4:DA:C8	2.54	0.42
1:A:48:ILE:HA	1:A:49:PRO:HD3	1.89	0.42
2:C:771:GLU:O	2:C:775:ARG:HB2	2.19	0.42
3:D:472:ALA:O	3:D:476:GLU:HG2	2.20	0.42
1:B:91:ASN:OD1	1:B:93:SER:HB2	2.20	0.42
2:C:195:LEU:O	2:C:199:VAL:HG23	2.20	0.42
2:C:571:LEU:HB2	2:C:574:ALA:HB2	2.02	0.42
3:D:1053[B]:PHE:CE2	3:D:1055:VAL:HB	2.54	0.42
5:F:96:LEU:O	5:F:100:VAL:HG23	2.20	0.42
3:D:636:GLN:CD	3:D:636:GLN:H	2.28	0.41
3:D:236:TYR:CE1	3:D:242:LEU:HD12	2.55	0.41
3:D:648:MET:HE2	3:D:648:MET:HB3	1.97	0.41
5:F:285:GLU:HA	5:F:286:PRO:HD3	1.84	0.41
2:C:397:GLU:N	2:C:633:GLN:OE1	2.46	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:192:ALA:HB1	3:D:193:PRO:HD2	2.02	0.41
5:F:105:LYS:HD3	5:F:179:GLU:HG2	2.01	0.41
6:G:15:DT:H2'	6:G:16:DC:C6	2.56	0.41
1:B:85:LEU:HG	1:B:87:VAL:HG23	2.02	0.41
2:C:94:LEU:HD22	2:C:118:ILE:HD11	2.03	0.41
2:C:1076:VAL:HG22	3:D:752:SER:HB3	2.02	0.41
2:C:1100:GLN:HG3	3:D:9:ARG:HH21	1.85	0.41
3:D:58:CYS:SG	3:D:62:LYS:N	2.93	0.41
3:D:313:MET:HE3	3:D:313:MET:HB3	1.90	0.41
3:D:1283:ILE:H	3:D:1283:ILE:HG13	1.60	0.41
5:F:208:SER:HB3	5:F:211:ASP:OD2	2.20	0.41
1:A:213:GLN:O	1:A:217:ILE:HG13	2.21	0.41
2:C:86:LYS:HB2	2:C:88:LEU:HG	2.02	0.41
2:C:198:ARG:HE	2:C:227:PHE:HA	1.85	0.41
3:D:65:ARG:HD3	3:D:65:ARG:HA	1.87	0.41
3:D:684:LYS:HE2	3:D:684:LYS:HB3	1.88	0.41
3:D:1267:ARG:HA	3:D:1268:PRO:HD3	1.85	0.41
5:F:144:ILE:HB	5:F:147:LEU:HD13	2.01	0.41
2:C:348:LEU:HD12	2:C:348:LEU:HA	1.93	0.41
2:C:413:LEU:HD21	2:C:451:LEU:HD13	2.01	0.41
2:C:971:LYS:HG2	2:C:988:VAL:HG22	2.02	0.41
3:D:1271:LYS:HD2	3:D:1331:ASP:HB2	2.03	0.41
3:D:38:LYS:HD3	3:D:38:LYS:HA	1.95	0.41
3:D:1084:THR:O	3:D:1088:THR:HG23	2.21	0.41
2:C:504:GLU:HG2	2:C:509:ALA:HB2	2.02	0.41
3:D:1189:ARG:HB3	3:D:1204:CYS:HA	2.03	0.41
4:E:37:ASN:OD1	4:E:37:ASN:N	2.37	0.41
2:C:76:PRO:HA	2:C:77:PRO:HD2	1.88	0.41
2:C:304:LEU:HB3	2:C:305:PRO:HD3	2.02	0.41
2:C:397:GLU:HB2	2:C:631:SER:HB2	2.02	0.41
2:C:834:GLN:HE22	3:D:724:GLN:NE2	2.18	0.41
2:C:976:ASP:HB3	2:C:979:THR:OG1	2.20	0.41
2:C:983:ILE:HG21	2:C:987:ILE:HD11	2.02	0.41
3:D:81:THR:OG1	3:D:82:LYS:N	2.53	0.41
3:D:904:VAL:HG13	3:D:905:PRO:O	2.21	0.41
3:D:1197:ARG:HB2	3:D:1398:TRP:CH2	2.56	0.41
1:A:64:GLU:HG2	1:A:76:VAL:HG22	2.02	0.41
2:C:157:ARG:HD3	2:C:157:ARG:HA	1.86	0.41
2:C:586:ARG:HH21	2:C:590:ASP:CG	2.29	0.41
2:C:670:GLN:HG2	2:C:699:PHE:CG	2.56	0.41
2:C:1044:GLY:O	3:D:1475:GLY:HA3	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:192:ALA:HB3	3:D:195:VAL:HB	2.03	0.41
3:D:750:PRO:HG2	3:D:756:GLN:NE2	2.36	0.41
3:D:796:ARG:HG3	3:D:861:GLN:HB3	2.03	0.41
3:D:803:GLY:HA2	3:D:827:ILE:HA	2.03	0.41
1:A:70:GLY:H	2:C:607:ASP:CG	2.29	0.40
1:B:141:GLU:OE1	1:B:161:ARG:NH2	2.54	0.40
2:C:302:VAL:O	2:C:305:PRO:HD2	2.20	0.40
2:C:1031:ARG:HA	3:D:622:ARG:HA	2.03	0.40
3:D:115:LEU:HD23	3:D:115:LEU:HA	1.83	0.40
3:D:658:LEU:HA	3:D:661:MET:HE3	2.03	0.40
3:D:1296:SER:HB3	3:D:1299:PHE:HB2	2.03	0.40
1:A:172:SER:HA	1:A:173:PRO:HD2	1.85	0.40
2:C:418:LEU:HD11	2:C:427:VAL:HG11	2.03	0.40
2:C:540:PHE:HB2	2:C:545:ASN:ND2	2.35	0.40
1:A:155:LYS:HA	1:A:155:LYS:HD2	1.87	0.40
2:C:223:ASP:OD1	2:C:225:SER:OG	2.36	0.40
2:C:674:VAL:HG12	2:C:869:VAL:HB	2.03	0.40
2:C:709:GLU:OE2	2:C:824:ARG:NH1	2.54	0.40
2:C:936:VAL:HG11	2:C:959:PRO:HB2	2.03	0.40
3:D:137:PRO:HA	3:D:452:ILE:HG23	2.03	0.40
5:F:364:ARG:HG3	5:F:390:PHE:CE2	2.57	0.40
1:A:102:LYS:HG3	1:A:139[B]:ASN:OD1	2.22	0.40
2:C:163:ILE:HG23	2:C:171:TRP:NE1	2.37	0.40
2:C:286:SER:OG	2:C:301:GLU:OE2	2.31	0.40
2:C:299:LYS:HE3	2:C:299:LYS:HB2	1.85	0.40
3:D:200:ASP:O	3:D:397:LYS:HG2	2.22	0.40
3:D:475:LYS:O	3:D:479:GLU:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	230/315 (73%)	225 (98%)	5 (2%)	0	100	100
1	B	221/315 (70%)	217 (98%)	4 (2%)	0	100	100
2	C	1111/1119 (99%)	1078 (97%)	33 (3%)	0	100	100
3	D	1485/1524 (97%)	1447 (97%)	38 (3%)	0	100	100
4	E	92/99 (93%)	89 (97%)	3 (3%)	0	100	100
5	F	344/443 (78%)	338 (98%)	6 (2%)	0	100	100
All	All	3483/3815 (91%)	3394 (97%)	89 (3%)	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	201/273 (74%)	197 (98%)	4 (2%)	48	72
1	B	196/273 (72%)	192 (98%)	4 (2%)	48	72
2	C	939/941 (100%)	903 (96%)	36 (4%)	29	62
3	D	1255/1279 (98%)	1221 (97%)	34 (3%)	39	68
4	E	83/88 (94%)	80 (96%)	3 (4%)	31	63
5	F	301/388 (78%)	292 (97%)	9 (3%)	36	66
All	All	2975/3242 (92%)	2885 (97%)	90 (3%)	36	66

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	66	SER
1	A	96	THR
1	A	134	GLU
1	B	7	LYS
1	B	10	VAL
1	B	34	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	154	GLU
2	C	11	GLU
2	C	138	SER
2	C	182	VAL
2	C	200	LEU
2	C	210	GLU
2	C	216	GLU
2	C	217	LEU
2	C	230	ARG
2	C	261	ILE
2	C	276	LYS
2	C	284	ARG
2	C	299	LYS
2	C	336	VAL
2	C	342	ASP
2	C	360	LEU
2	C	384	GLU
2	C	390	GLN
2	C	409	ARG
2	C	421	GLU
2	C	454	SER
2	C	513	VAL
2	C	575	GLN
2	C	595	LEU
2	C	617	ASP
2	C	620	LEU
2	C	657	ASP
2	C	670	GLN
2	C	674	VAL
2	C	722	ILE
2	C	775	ARG
2	C	807[A]	ARG
2	C	807[B]	ARG
2	C	888	THR
2	C	952	LEU
2	C	974	LEU
2	C	1076	VAL
3	D	35	ARG
3	D	71	LYS
3	D	106	LYS
3	D	231	VAL
3	D	354	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	D	407	VAL
3	D	531	ASP
3	D	632	VAL
3	D	636	GLN
3	D	687	VAL
3	D	709	HIS
3	D	754	PHE
3	D	778	LEU
3	D	784	ASP
3	D	810	GLU
3	D	829	VAL
3	D	894	LYS
3	D	904	VAL
3	D	907	GLU
3	D	956	ILE
3	D	986	ARG
3	D	1067	VAL
3	D	1129	THR
3	D	1184	GLN
3	D	1188	VAL
3	D	1223	ILE
3	D	1252	ILE
3	D	1253	THR
3	D	1267	ARG
3	D	1283	ILE
3	D	1290	LEU
3	D	1295	GLU
3	D	1307	LYS
3	D	1408	ILE
4	E	37	ASN
4	E	49	GLN
4	E	50	THR
5	F	88	ILE
5	F	120	THR
5	F	162	LYS
5	F	295	MET
5	F	310	ILE
5	F	321	ILE
5	F	325	LYS
5	F	380	GLU
5	F	399	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16)

such sidechains are listed below:

Mol	Chain	Res	Type
2	C	187	ASN
2	C	219	GLN
2	C	556	ASN
3	D	348	GLN
3	D	350	HIS
3	D	611	GLN
3	D	714	GLN
3	D	768	ASN
3	D	976	GLN
3	D	994	GLN
3	D	1124	GLN
3	D	1184	GLN
3	D	1195	GLN
3	D	1227	GLN
5	F	83	GLN
5	F	218	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
8	I	3/5 (60%)	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 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%)	0.13	1 (0%) 88 79	26, 50, 76, 114	1 (0%)
1	B	223/315 (70%)	0.15	3 (1%) 75 55	24, 53, 89, 105	0
2	C	1112/1119 (99%)	-0.08	12 (1%) 78 61	7, 42, 100, 146	3 (0%)
3	D	1486/1524 (97%)	-0.03	24 (1%) 70 51	5, 38, 102, 136	4 (0%)
4	E	94/99 (94%)	-0.01	1 (1%) 78 61	16, 44, 84, 93	0
5	F	346/443 (78%)	0.47	35 (10%) 12 8	20, 59, 129, 145	0
6	G	17/21 (80%)	0.00	1 (5%) 28 18	9, 29, 134, 134	0
7	H	24/27 (88%)	0.39	1 (4%) 40 25	43, 64, 106, 139	0
8	I	4/5 (80%)	-1.36	0 100 100	10, 13, 18, 23	0
All	All	3537/3868 (91%)	0.03	78 (2%) 62 42	5, 46, 105, 146	8 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	184	GLU	4.3
2	C	102[A]	HIS	3.9
3	D	144	GLY	3.9
2	C	64	LEU	3.8
3	D	1129	THR	3.5
5	F	373	LYS	3.4
5	F	396	ARG	3.4
7	H	24	DC	3.4
5	F	404	ALA	3.3
5	F	389	PHE	3.2
5	F	386	VAL	3.2
5	F	399	GLN	3.1
2	C	63	GLY	3.1
6	G	4	DG	3.0
5	F	414	ARG	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
3	D	420	VAL	3.0
5	F	141	VAL	3.0
5	F	369	LEU	2.9
2	C	769	PRO	2.9
5	F	388	ALA	2.9
3	D	434	ARG	2.8
5	F	324	GLU	2.8
5	F	393	THR	2.8
5	F	395	GLU	2.8
5	F	375	LEU	2.8
1	B	228	PRO	2.7
3	D	1313	VAL	2.7
5	F	397	ILE	2.7
5	F	415	THR	2.6
3	D	235	ALA	2.6
4	E	95	VAL	2.6
2	C	103	LYS	2.6
5	F	392	VAL	2.6
5	F	142	ARG	2.5
2	C	908	GLY	2.4
2	C	101	ILE	2.4
3	D	1283	ILE	2.4
2	C	773	LEU	2.4
5	F	376	ILE	2.4
3	D	415	VAL	2.4
5	F	139	ALA	2.3
2	C	56	GLU	2.3
3	D	241	ILE	2.3
3	D	431	VAL	2.3
3	D	172	PRO	2.3
3	D	531	ASP	2.3
5	F	400	ILE	2.3
5	F	413	SER	2.3
1	B	6	LEU	2.2
5	F	358	LEU	2.2
2	C	188	LYS	2.2
2	C	275	TYR	2.2
5	F	390	PHE	2.2
5	F	410	TYR	2.2
5	F	166	LEU	2.2
1	B	187	GLY	2.2
5	F	153	PRO	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
5	F	391	GLY	2.2
5	F	405	LEU	2.2
3	D	183	GLU	2.2
3	D	407	VAL	2.1
5	F	322	GLY	2.1
3	D	1252	ILE	2.1
2	C	364	GLU	2.1
3	D	173	PRO	2.1
5	F	378	GLY	2.1
5	F	145	PRO	2.1
3	D	1299	PHE	2.1
3	D	429	SER	2.1
3	D	175	VAL	2.1
3	D	978	TYR	2.1
1	A	139[A]	ASN	2.0
3	D	310	LEU	2.0
5	F	383	LEU	2.0
3	D	316	GLN	2.0
3	D	203	ALA	2.0
5	F	403	LYS	2.0
5	F	144	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	0.83	0.18	54,54,54,54	0
9	MG	F	2001	1/1	0.88	0.07	48,48,48,48	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	MG	D	2004	1/1	0.96	0.18	28,28,28,28	0
9	MG	D	2003	1/1	0.98	0.04	8,8,8,8	0
10	ZN	D	2002	1/1	0.99	0.03	103,103,103,103	0
10	ZN	D	2001	1/1	1.00	0.02	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.