



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 01:24 AM BST

PDB ID : 6L74
Title : Thermus thermophilus initial transcription complex comprising sigma A and 5'-triphosphate RNA of 2 nt
Authors : Zhang, Y.; Li, L.; Ebright, R.H.
Deposited on : 2019-10-31
Resolution : 3.12 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

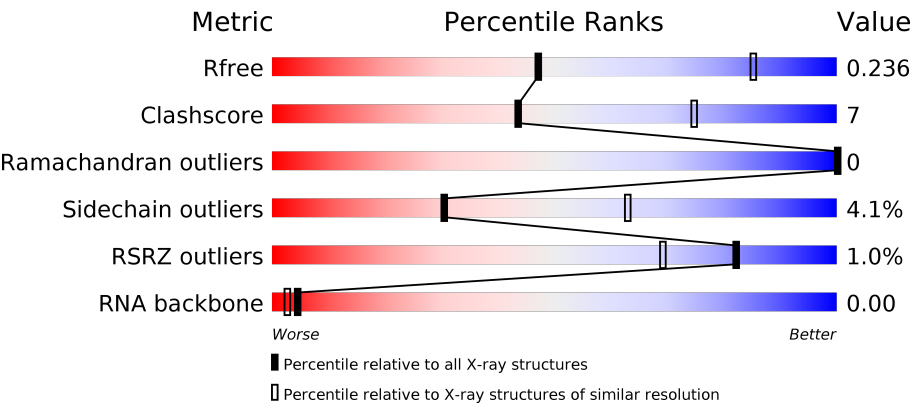
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.12 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.







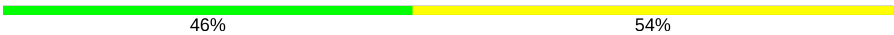
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	130704	1292 (3.14-3.10)
Clashscore	141614	1389 (3.14-3.10)
Ramachandran outliers	138981	1337 (3.14-3.10)
Sidechain outliers	138945	1337 (3.14-3.10)
RSRZ outliers	127900	1260 (3.14-3.10)
RNA backbone	3102	1134 (3.44-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	315	<div><div>%</div><div><div></div><div>55%</div><div>17%</div><div>.</div><div>27%</div></div></div>
1	B	315	<div><div>%</div><div><div></div><div>57%</div><div>14%</div><div>.</div><div>28%</div></div></div>
2	C	1119	<div><div>%</div><div><div></div><div>79%</div><div>19%</div><div>..</div></div></div>
3	D	1524	<div><div>%</div><div><div></div><div>79%</div><div>18%</div><div>..</div></div></div>

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Mol	Chain	Length	Quality of chain
4	E	99	 87% 9% .
5	F	443	 % 63% 14% • 22%
6	G	16	 56% 44%
7	I	2	 100%
8	H	24	 46% 54%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	MG	B	2002	-	-	-	X

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 28744 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
			1789	1143	310	334	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
			8774	5550	1565	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	1	0
			11729	7435	2066	2193	35			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	95	Total	C	N	O	S	0	0	0
			770	491	133	142	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(*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*CP*GP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	16	Total	C	N	O	P	0	0	0
			328	156	63	94	15			

- Molecule 7 is a RNA chain called RNA (5'-D(*(GTP))-R(P*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	I	2	Total	C	N	O	P	0	0	0
			54	20	10	20	4			

- Molecule 8 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*C)-3').

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

- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	2	Total Mg 2 2	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 (three-letter code: ZN) (formula: Zn).

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

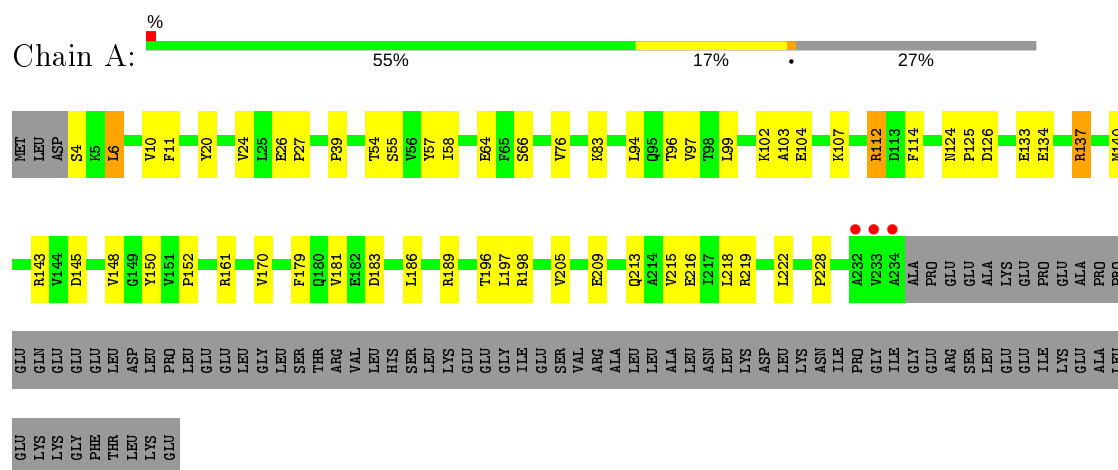
- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	11	Total O 11 11	0	0
11	B	10	Total O 10 10	0	0
11	C	55	Total O 55 55	0	0
11	D	80	Total O 80 80	0	0
11	E	5	Total O 5 5	0	0
11	F	10	Total O 10 10	0	0
11	G	6	Total O 6 6	0	0
11	I	4	Total O 4 4	0	0
11	H	1	Total O 1 1	0	0

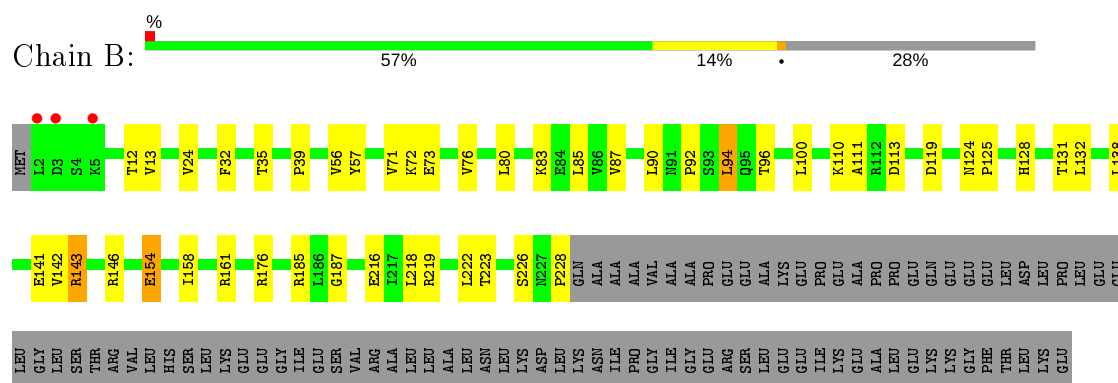
3 Residue-property plots

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.

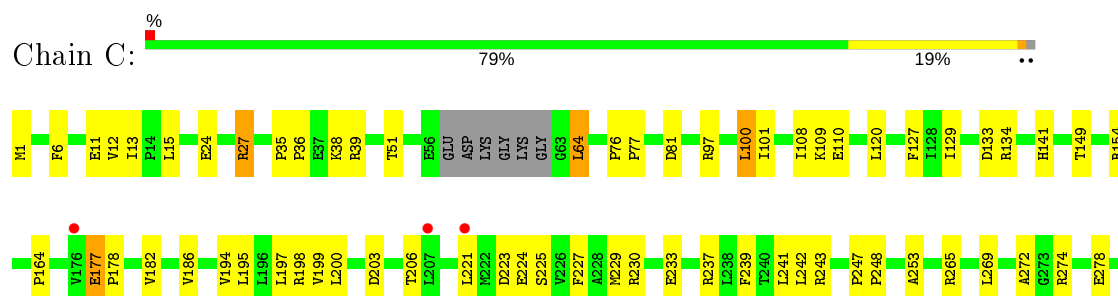
- Molecule 1: DNA-directed RNA polymerase subunit alpha

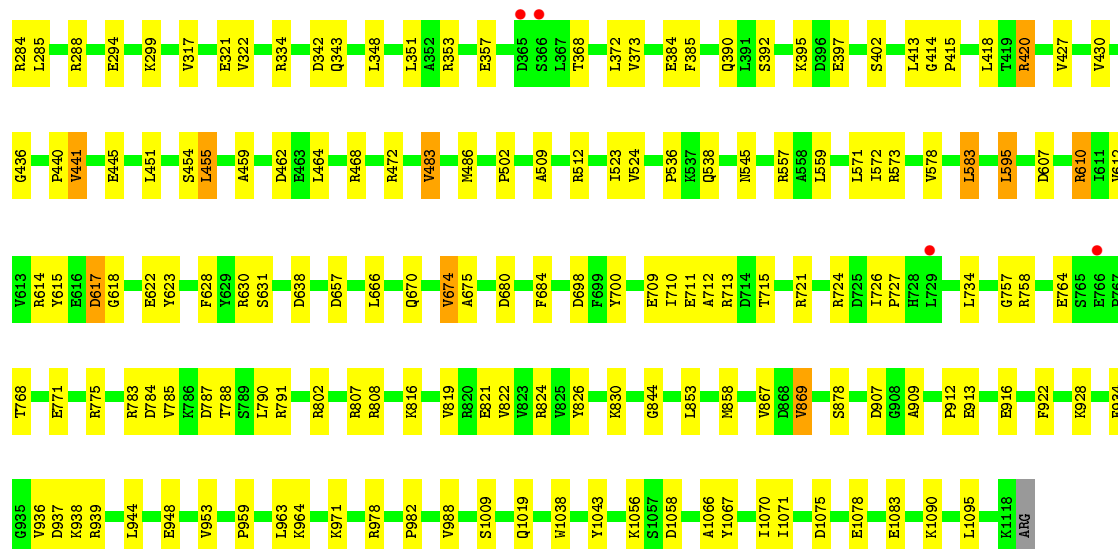


- Molecule 1: DNA-directed RNA polymerase subunit alpha

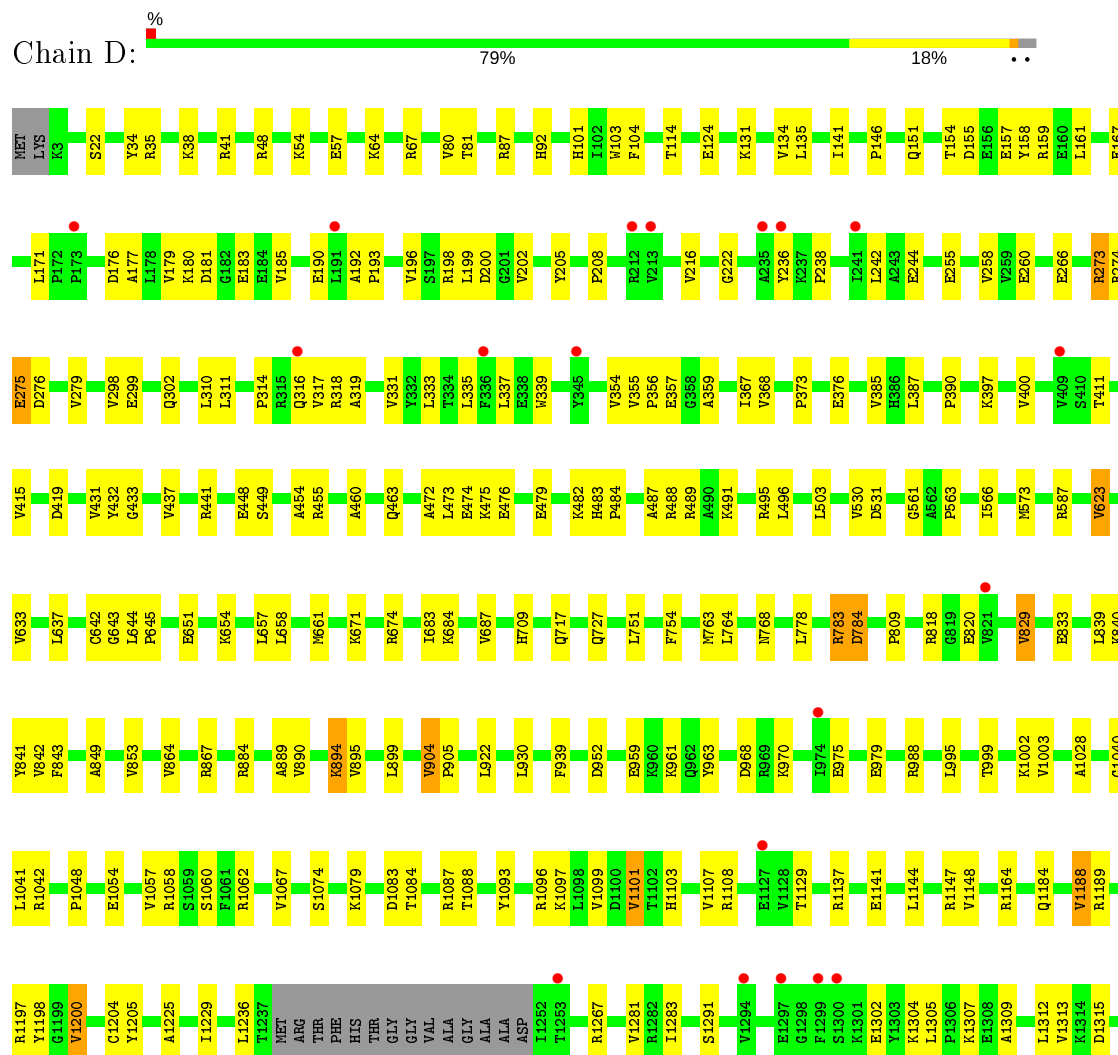


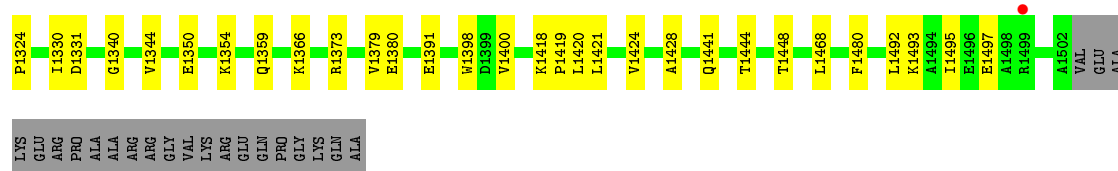
- Molecule 2: DNA-directed RNA polymerase subunit beta





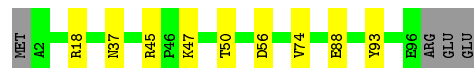
• Molecule 3: DNA-directed RNA polymerase subunit beta'





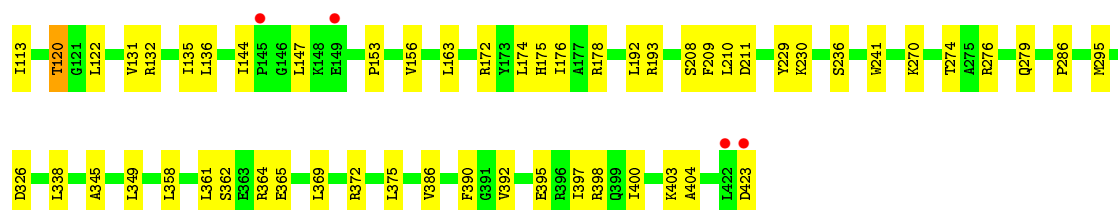
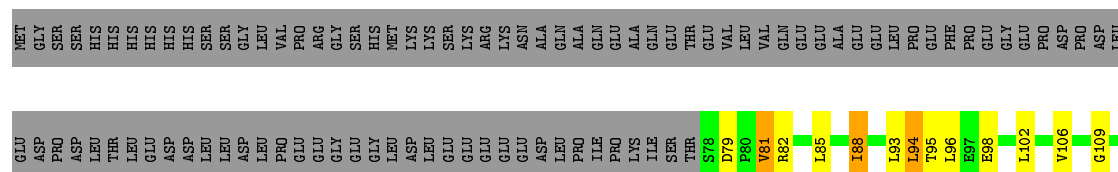
- Molecule 4: DNA-directed RNA polymerase subunit omega

Chain E: 87% 9%



- Molecule 5: RNA polymerase sigma factor SigA

Chain F: 63% 14% 22%



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

Chain G: 56% 44%



- Molecule 7: RNA (5'-D(*(GTP))-R(P*A)-3')

Chain I: 100%

There are no outlier residues recorded for this chain.

- Molecule 8: 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*C)-3')

Chain H: 46% 54%

T1	A2	T3	A4	A5	G8	G9	T15	C16	A17	C18	G19	G20	A21	C24
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	183.07Å 103.50Å 295.13Å 90.00° 99.17° 90.00°	Depositor
Resolution (Å)	44.92 – 3.12 44.92 – 3.12	Depositor EDS
% Data completeness (in resolution range)	99.3 (44.92-3.12) 99.3 (44.92-3.12)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.191 , 0.236 0.191 , 0.236	Depositor DCC
R_{free} test set	2128 reflections (2.20%)	wwPDB-VP
Wilson B-factor (Å ²)	80.1	Xtriage
Anisotropy	0.551	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for 1/2*h-3/2*k,-1/2*h-1/2*k,-1/2*h +1/2*k-l 0.017 for 1/2*h+3/2*k,1/2*h-1/2*k,-1/2*h- 1/2*k-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	28744	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: GTP, 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	0.26	0/1841	0.47	0/2504
1	B	0.26	0/1821	0.45	0/2476
2	C	0.26	0/8941	0.45	0/12092
3	D	0.26	0/11938	0.45	0/16143
4	E	0.25	0/784	0.41	0/1057
5	F	0.25	0/2852	0.41	0/3837
6	G	0.57	0/368	0.83	0/567
7	I	0.28	0/24	0.74	0/35
8	H	0.57	0/556	0.94	0/858
All	All	0.27	0/29125	0.47	0/39569

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	829	VAL	Peptide

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	34	0
1	B	1789	0	1841	35	0
2	C	8774	0	8877	121	0
3	D	11729	0	11949	167	0
4	E	770	0	784	6	0
5	F	2807	0	2882	41	0
6	G	328	0	181	10	0
7	I	54	0	22	0	0
8	H	495	0	272	19	0
9	B	2	0	0	0	0
9	D	2	0	0	0	0
9	F	1	0	0	0	0
10	D	2	0	0	0	0
11	A	11	0	0	0	0
11	B	10	0	0	0	0
11	C	55	0	0	2	0
11	D	80	0	0	5	0
11	E	5	0	0	0	0
11	F	10	0	0	0	0
11	G	6	0	0	0	0
11	H	1	0	0	0	0
11	I	4	0	0	0	0
All	All	28744	0	28671	383	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 (383) 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:7:DT:H3	8:H:21:DA:H61	1.16	0.94
6:G:10:DG:H1	8:H:18:DC:H42	1.18	0.91
8:H:16:DC:H2'	8:H:17:DA:C8	2.21	0.75
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.70	0.73
3:D:135:LEU:HD13	3:D:463:GLN:HG2	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:361:LEU:HB3	5:F:365:GLU:HG3	1.72	0.72
8:H:18:DC:H2'	8:H:19:DG:C8	2.25	0.71
1:B:132:LEU:HD21	1:B:138:LEU:HB2	1.73	0.71
3:D:1236:LEU:HA	3:D:1359:GLN:HG3	1.72	0.71
3:D:134:VAL:HG22	3:D:151:GLN:H	1.55	0.70
2:C:628:PHE:H	2:C:638:ASP:HB3	1.57	0.69
6:G:15:DT:H2'	6:G:16:DC:C6	2.28	0.69
5:F:397:ILE:HD12	5:F:400:ILE:HD11	1.74	0.68
3:D:1048:PRO:O	3:D:1079:LYS:NZ	2.25	0.68
2:C:194:VAL:HG22	2:C:221:LEU:HD12	1.74	0.67
2:C:1083:GLU:OE2	3:D:87:ARG:NH2	2.28	0.67
2:C:713:ARG:HA	2:C:819:VAL:HA	1.77	0.66
3:D:255:GLU:OE1	3:D:274:ARG:NH2	2.23	0.66
2:C:709:GLU:OE2	2:C:824:ARG:NH1	2.29	0.66
3:D:489:ARG:NH1	3:D:1391:GLU:OE2	2.29	0.66
1:B:94:LEU:O	1:B:146:ARG:NH2	2.28	0.65
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.77	0.65
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.29	0.65
5:F:95:THR:HB	5:F:98:GLU:HG3	1.78	0.65
2:C:24:GLU:OE2	2:C:27:ARG:NH2	2.31	0.64
2:C:462:ASP:HB3	2:C:468:ARG:HD2	1.78	0.64
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.79	0.64
2:C:617:ASP:OD1	2:C:617:ASP:N	2.28	0.64
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.80	0.64
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.31	0.64
3:D:1281:VAL:HG21	3:D:1313:VAL:HG21	1.79	0.64
2:C:802:ARG:HB2	2:C:826:TYR:HB2	1.80	0.63
2:C:680:ASP:OD2	2:C:978:ARG:NH2	2.31	0.63
8:H:19:DG:H2''	8:H:20:DG:H2'	1.81	0.62
1:A:112:ARG:HG3	1:A:125:PRO:HB2	1.82	0.62
4:E:45:ARG:NH1	4:E:56:ASP:OD2	2.33	0.62
2:C:334:ARG:NH2	2:C:342:ASP:OD2	2.33	0.61
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.81	0.61
3:D:411:THR:O	5:F:178:ARG:NH1	2.30	0.61
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.83	0.61
2:C:1056:LYS:HB2	3:D:623:VAL:HG22	1.83	0.60
8:H:18:DC:H2''	8:H:19:DG:H5'	1.83	0.60
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.83	0.60
1:A:104:GLU:HB3	1:A:137:ARG:HG3	1.82	0.60
2:C:711:GLU:HG2	2:C:822:VAL:HG22	1.83	0.60
5:F:400:ILE:HA	5:F:403:LYS:HG2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:657:LEU:HG	3:D:661:MET:HE2	1.84	0.60
2:C:353:ARG:NH1	2:C:357:GLU:OE2	2.36	0.59
3:D:1042:ARG:HB3	3:D:1057:VAL:HB	1.85	0.59
3:D:1101:VAL:HG21	3:D:1424:VAL:HG22	1.85	0.59
2:C:787:ASP:OD2	2:C:791:ARG:NH2	2.35	0.59
1:A:216:GLU:OE2	1:A:219:ARG:NH2	2.36	0.59
3:D:1083:ASP:OD1	3:D:1087:ARG:NH2	2.36	0.59
2:C:35:PRO:HG2	2:C:38:LYS:HD2	1.85	0.58
3:D:155:ASP:OD1	3:D:159:ARG:NH2	2.36	0.58
4:E:47:LYS:NZ	4:E:56:ASP:OD1	2.34	0.58
5:F:94:LEU:HD21	5:F:102:LEU:HD12	1.85	0.58
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.86	0.58
2:C:64:LEU:HG	2:C:100:LEU:HD21	1.83	0.58
3:D:1324:PRO:HG3	3:D:1330:ILE:HD11	1.85	0.57
1:A:58:ILE:HG12	1:A:140:MET:HG2	1.85	0.57
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.86	0.57
5:F:365:GLU:HB2	5:F:404:ALA:HB2	1.85	0.57
2:C:768:THR:OG1	2:C:771:GLU:OE1	2.21	0.56
2:C:12:VAL:HG21	2:C:472:ARG:HD3	1.87	0.56
1:B:92:PRO:O	1:B:146:ARG:NH2	2.38	0.56
3:D:316:GLN:NE2	11:D:2108:HOH:O	2.39	0.56
2:C:413:LEU:HD21	2:C:451:LEU:HD13	1.87	0.56
6:G:15:DT:H2'	6:G:16:DC:H6	1.71	0.56
3:D:433:GLY:HA2	3:D:449:SER:H	1.72	0.55
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.88	0.55
1:A:209:GLU:O	1:A:213:GLN:HG2	2.06	0.55
2:C:164:PRO:HA	2:C:269:LEU:HD23	1.88	0.55
2:C:614:ARG:NH2	2:C:618:GLY:O	2.39	0.55
3:D:1283:ILE:HG12	3:D:1315:ASP:HB2	1.88	0.55
2:C:1019:GLN:HG2	2:C:1058:ASP:HB3	1.88	0.55
2:C:607:ASP:HB2	2:C:610:ARG:NH1	2.20	0.55
1:A:103:ALA:HB1	1:A:107:LYS:HD3	1.88	0.55
1:B:216:GLU:OE1	1:B:219:ARG:NH2	2.37	0.55
2:C:197:LEU:HD12	2:C:221:LEU:HD11	1.89	0.55
3:D:487:ALA:O	3:D:491:LYS:HG2	2.07	0.55
3:D:658:LEU:HA	3:D:661:MET:HE3	1.90	0.55
3:D:683:ILE:HG23	3:D:687:VAL:HG11	1.89	0.54
3:D:134:VAL:HG12	3:D:454:ALA:HB2	1.89	0.54
3:D:479:GLU:HA	3:D:482:LYS:HE2	1.90	0.54
1:A:39:PRO:HG3	1:B:39:PRO:HG3	1.88	0.54
5:F:93:LEU:HD21	5:F:193:ARG:HD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:224:GLU:CD	2:C:224:GLU:H	2.11	0.54
3:D:970:LYS:HD3	3:D:995:LEU:HD13	1.90	0.54
8:H:15:DT:H2"	8:H:16:DC:C6	2.42	0.54
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.89	0.54
1:B:80:LEU:HD21	3:D:842:VAL:HG12	1.89	0.54
3:D:1093:TYR:OH	3:D:1441:GLN:OE1	2.20	0.54
3:D:895:VAL:HG11	3:D:922:LEU:HD21	1.89	0.54
6:G:12:DG:N2	8:H:16:DC:N3	2.50	0.54
2:C:97:ARG:NH1	2:C:110:GLU:OE1	2.41	0.53
3:D:176:ASP:OD1	3:D:177:ALA:N	2.39	0.53
3:D:561:GLY:HA3	5:F:132:ARG:HD3	1.90	0.53
3:D:57:GLU:HG3	3:D:64:LYS:HG2	1.89	0.53
2:C:572:ILE:HG13	2:C:573:ARG:HG3	1.89	0.53
3:D:1096:ARG:NH1	11:D:2103:HOH:O	2.33	0.53
2:C:684:PHE:HB3	3:D:633:VAL:HG21	1.90	0.53
2:C:272:ALA:HA	2:C:464:LEU:HD13	1.89	0.53
1:A:215:VAL:HG13	1:B:222:LEU:HD22	1.91	0.53
2:C:294:GLU:HB3	2:C:299:LYS:HE2	1.91	0.53
2:C:545:ASN:HB3	2:C:583:LEU:HD12	1.91	0.53
5:F:270:LYS:O	5:F:274:THR:OG1	2.25	0.52
5:F:88:ILE:HG23	5:F:193:ARG:HG2	1.92	0.52
3:D:473:LEU:HD21	3:D:495:ARG:HH21	1.74	0.52
2:C:15:LEU:HD12	2:C:583:LEU:HD21	1.92	0.52
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.92	0.52
5:F:135:ILE:HD11	5:F:178:ARG:HB3	1.91	0.52
2:C:536:PRO:HB3	3:D:1067:VAL:HG21	1.92	0.52
2:C:1043:TYR:CG	3:D:763:MET:HG2	2.45	0.52
6:G:12:DG:H1	8:H:16:DC:H42	1.57	0.52
3:D:274:ARG:NH2	3:D:279:VAL:HG21	2.24	0.52
2:C:1067:TYR:OH	3:D:674[A]:ARG:NH1	2.43	0.52
3:D:573:MET:SD	5:F:210:LEU:HB3	2.50	0.52
2:C:1071:ILE:HD11	5:F:345:ALA:HB1	1.92	0.51
3:D:208:PRO:HA	3:D:390:PRO:HA	1.92	0.51
2:C:1056:LYS:HE2	3:D:751:LEU:HG	1.92	0.51
3:D:472:ALA:O	3:D:476:GLU:HG2	2.10	0.51
2:C:721:ARG:HH22	2:C:785:VAL:HG11	1.74	0.51
3:D:356:PRO:HB3	3:D:441:ARG:HA	1.91	0.51
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.93	0.51
2:C:948:GLU:HG3	2:C:953:VAL:HG23	1.93	0.51
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.11	0.51
3:D:114:THR:HG23	3:D:495:ARG:HG2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1038:TRP:CE2	3:D:1099:VAL:HG11	2.46	0.50
2:C:198:ARG:HE	2:C:227:PHE:HA	1.75	0.50
1:A:20:TYR:OH	1:A:198:ARG:HD2	2.11	0.50
5:F:153:PRO:HA	5:F:156:VAL:HG22	1.92	0.50
5:F:163:LEU:HD13	5:F:174:LEU:HD13	1.92	0.50
2:C:1009:SER:HB3	3:D:651:GLU:O	2.11	0.50
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.92	0.50
2:C:724:ARG:NH2	2:C:734:LEU:O	2.44	0.50
2:C:758:ARG:HH21	2:C:788:THR:HB	1.76	0.50
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.94	0.49
1:B:176:ARG:HD3	3:D:884:ARG:NH2	2.27	0.49
3:D:784:ASP:HB2	3:D:939:PHE:CE2	2.47	0.49
4:E:37:ASN:N	4:E:37:ASN:OD1	2.33	0.49
2:C:628:PHE:H	2:C:638:ASP:CB	2.22	0.49
3:D:1137:ARG:O	3:D:1141:GLU:HG3	2.12	0.49
1:B:154:GLU:HG3	3:D:840:LYS:NZ	2.26	0.49
1:A:4:SER:O	1:A:189:ARG:NH2	2.38	0.49
2:C:101:ILE:HG12	2:C:108:ILE:HG12	1.94	0.49
2:C:134:ARG:NH1	2:C:392:SER:O	2.45	0.49
2:C:937:ASP:OD1	2:C:938:LYS:N	2.46	0.49
5:F:241:TRP:CE2	8:H:1:DT:H4'	2.48	0.49
1:B:223:THR:O	1:B:226:SER:OG	2.26	0.48
2:C:274:ARG:NH2	2:C:285:LEU:O	2.46	0.48
3:D:1097:LYS:O	3:D:1101:VAL:HG22	2.13	0.48
3:D:637:LEU:HD13	3:D:642:CYS:HA	1.95	0.48
8:H:20:DG:H1'	8:H:21:DA:C8	2.48	0.48
2:C:944:LEU:HD21	2:C:963:LEU:HD23	1.95	0.48
3:D:1380:GLU:HB2	3:D:1420:LEU:HD22	1.95	0.48
6:G:10:DG:H1	8:H:18:DC:N4	1.99	0.48
2:C:615:TYR:OH	2:C:623:TYR:OH	2.22	0.48
5:F:79:ASP:OD2	8:H:8:DG:N1	2.33	0.48
2:C:274:ARG:HH12	2:C:284:ARG:HH22	1.61	0.47
3:D:244:GLU:HA	3:D:310:LEU:HA	1.96	0.47
1:B:143:ARG:NH1	1:B:158:ILE:HD12	2.29	0.47
2:C:351:LEU:HD11	2:C:373:VAL:HG13	1.97	0.47
3:D:158:TYR:CE1	3:D:454:ALA:HB3	2.49	0.47
3:D:415:VAL:HG13	3:D:419:ASP:HB2	1.95	0.47
2:C:133:ASP:HB3	2:C:395:LYS:HD2	1.96	0.47
2:C:939:ARG:HG2	2:C:982:PRO:HD3	1.96	0.47
3:D:483:HIS:CG	3:D:484:PRO:HD2	2.50	0.47
5:F:144:ILE:HB	5:F:147:LEU:HD13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:GLU:OE1	1:B:161:ARG:NH2	2.47	0.47
1:B:32:PHE:HA	1:B:35:THR:HB	1.97	0.47
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.76	0.47
2:C:455:LEU:HD22	2:C:459:ALA:HB3	1.97	0.47
3:D:890:VAL:HB	3:D:922:LEU:HD13	1.96	0.47
3:D:1084:THR:O	3:D:1088:THR:HG23	2.15	0.47
3:D:1493:LYS:O	3:D:1497:GLU:HG2	2.14	0.47
3:D:1480:PHE:O	4:E:18:ARG:NH2	2.48	0.47
5:F:172:ARG:O	5:F:176:ILE:HG12	2.14	0.47
5:F:208:SER:HB3	5:F:211:ASP:OD2	2.14	0.47
2:C:76:PRO:HG3	2:C:120:LEU:HD12	1.97	0.47
2:C:129:ILE:HB	2:C:134:ARG:HD2	1.97	0.47
3:D:190:GLU:HA	3:D:196:VAL:HA	1.97	0.47
8:H:20:DG:H3'	8:H:20:DG:OP2	2.14	0.47
3:D:975:GLU:O	3:D:979:GLU:HG2	2.15	0.46
5:F:395:GLU:OE2	5:F:398:ARG:NH2	2.48	0.46
2:C:928:LYS:HB2	2:C:928:LYS:HE3	1.76	0.46
3:D:1189:ARG:HB3	3:D:1204:CYS:HA	1.97	0.46
3:D:200:ASP:O	3:D:397:LYS:HG2	2.15	0.46
1:B:83:LYS:NZ	3:D:842:VAL:O	2.48	0.46
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.50	0.46
1:B:12:THR:HB	1:B:24:VAL:HB	1.97	0.46
2:C:440:PRO:HB2	3:D:1074:SER:OG	2.15	0.46
3:D:1302:GLU:OE1	3:D:1304:LYS:HE3	2.15	0.46
3:D:783:ARG:HD2	3:D:1028:ALA:O	2.14	0.46
2:C:36:PRO:HA	2:C:39:ARG:HG3	1.98	0.46
2:C:674:VAL:HG22	2:C:869:VAL:HG22	1.96	0.46
3:D:236:TYR:CE1	3:D:242:LEU:HD12	2.50	0.46
5:F:270:LYS:HG2	5:F:295:MET:HE1	1.98	0.46
1:A:228:PRO:HB3	1:B:13:VAL:HG21	1.97	0.46
1:A:83:LYS:NZ	2:C:698:ASP:OD1	2.48	0.46
5:F:362:SER:OG	5:F:365:GLU:HG2	2.16	0.46
3:D:258:VAL:HG12	3:D:273:ARG:O	2.16	0.46
3:D:432:TYR:O	3:D:448:GLU:HA	2.16	0.46
3:D:784:ASP:HB2	3:D:939:PHE:HE2	1.81	0.46
3:D:764:LEU:O	3:D:768:ASN:ND2	2.49	0.46
1:A:57:TYR:CD1	1:A:161:ARG:HD2	2.51	0.45
3:D:1267:ARG:NE	3:D:1331:ASP:OD2	2.50	0.45
3:D:298:VAL:HG12	3:D:302:GLN:NE2	2.31	0.45
1:A:133:GLU:HG2	1:A:134:GLU:H	1.81	0.45
1:A:64:GLU:OE2	2:C:830:LYS:NZ	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:278:GLU:OE2	2:C:284:ARG:NH2	2.48	0.45
3:D:809:PRO:HB3	3:D:839:LEU:HD13	1.98	0.45
3:D:1099:VAL:O	3:D:1103:HIS:HB3	2.17	0.45
3:D:141:ILE:HA	3:D:146:PRO:HA	1.98	0.45
3:D:644:LEU:HD12	3:D:645:PRO:HD2	1.97	0.45
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.98	0.45
2:C:922:PHE:CD2	2:C:964:LYS:HG3	2.52	0.45
3:D:437:VAL:HG11	5:F:175:HIS:CD2	2.52	0.45
3:D:843:PHE:HE1	3:D:864:VAL:HG21	1.81	0.45
2:C:13:ILE:HD13	2:C:483:VAL:HG21	1.99	0.45
2:C:721:ARG:HH12	2:C:785:VAL:HG11	1.82	0.45
3:D:1344:VAL:HG11	3:D:1421:LEU:HD22	1.98	0.45
2:C:816:LYS:O	2:C:819:VAL:HG13	2.17	0.45
3:D:1137:ARG:HB3	11:D:2125:HOH:O	2.17	0.45
2:C:243:ARG:NH1	8:H:9:DG:O6	2.50	0.45
3:D:988:ARG:NH2	3:D:1054:GLU:OE2	2.47	0.45
3:D:22:SER:HB2	3:D:92:HIS:HB3	1.97	0.45
3:D:894:LYS:HD2	3:D:894:LYS:H	1.81	0.45
2:C:436:GLY:HA2	2:C:538:GLN:O	2.17	0.45
2:C:807:ARG:HG2	2:C:821:GLU:HB3	1.98	0.45
8:H:3:DT:H2'	8:H:4:DA:C8	2.51	0.45
3:D:1197:ARG:HB2	3:D:1398:TRP:CH2	2.53	0.44
3:D:373:PRO:HA	3:D:376:GLU:HG3	1.98	0.44
3:D:455:ARG:HB2	3:D:460:ALA:HB2	2.00	0.44
2:C:971:LYS:HG2	2:C:988:VAL:HG22	1.99	0.44
3:D:179:VAL:O	3:D:205:TYR:OH	2.34	0.44
2:C:502:PRO:HB2	2:C:509:ALA:HB3	1.99	0.44
3:D:1444:THR:O	3:D:1448:THR:HG23	2.17	0.44
3:D:157:GLU:O	3:D:161:LEU:HG	2.18	0.44
2:C:727:PRO:HB3	2:C:783:ARG:HD3	2.00	0.44
3:D:684:LYS:O	3:D:687:VAL:HG12	2.18	0.44
1:A:10:VAL:HG12	1:A:26:GLU:O	2.18	0.44
1:A:124:ASN:OD1	1:A:124:ASN:N	2.50	0.44
1:A:24:VAL:HG22	1:A:196:THR:HG23	2.00	0.44
2:C:712:ALA:HB3	2:C:821:GLU:HG3	2.00	0.44
3:D:131:LYS:NZ	3:D:154:THR:HG22	2.32	0.44
5:F:279:GLN:HB3	5:F:286:PRO:HD3	1.99	0.44
1:B:76:VAL:O	1:B:80:LEU:HG	2.18	0.44
2:C:390:GLN:HG2	2:C:414:GLY:HA2	2.00	0.44
3:D:1373:ARG:HD3	11:D:2140:HOH:O	2.17	0.44
3:D:566:ILE:HD11	5:F:192:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:583:LEU:HA	2:C:583:LEU:HD23	1.73	0.44
3:D:1101:VAL:HG13	3:D:1428:ALA:HB2	1.98	0.44
1:B:110:LYS:HD3	1:B:128:HIS:HA	2.00	0.43
1:B:57:TYR:CD1	1:B:161:ARG:HD2	2.53	0.43
1:B:72:LYS:HG3	1:B:73:GLU:N	2.34	0.43
2:C:784:ASP:N	2:C:784:ASP:OD1	2.45	0.43
3:D:563:PRO:HD2	3:D:566:ILE:HD12	2.00	0.43
5:F:229:TYR:CZ	5:F:230:LYS:HD3	2.53	0.43
6:G:12:DG:H1	8:H:16:DC:N4	2.16	0.43
1:A:54:THR:HG21	1:A:145:ASP:HB2	1.99	0.43
5:F:82:ARG:HB2	8:H:8:DG:O6	2.18	0.43
1:A:150:TYR:CE2	1:A:152:PRO:HG3	2.54	0.43
1:B:100:LEU:HG	1:B:141:GLU:HG2	2.00	0.43
2:C:557:ARG:NH2	11:C:1205:HOH:O	2.34	0.43
1:A:181:VAL:HG12	2:C:938:LYS:HD2	2.00	0.43
3:D:208:PRO:HG3	3:D:387:LEU:HD22	1.99	0.43
3:D:889:ALA:HB1	3:D:930:LEU:HA	1.99	0.43
3:D:999:THR:O	3:D:1003:VAL:HG13	2.18	0.43
5:F:88:ILE:HD11	5:F:192:LEU:HD13	1.99	0.43
1:B:216:GLU:CD	1:B:219:ARG:HH21	2.21	0.43
2:C:154:ARG:NH1	2:C:178:PRO:HG3	2.33	0.43
3:D:904:VAL:HG22	3:D:905:PRO:HD2	2.00	0.43
3:D:671:LYS:NZ	5:F:423:ASP:OD1	2.38	0.43
3:D:1291:SER:OG	3:D:1304:LYS:HG2	2.18	0.43
3:D:1350:GLU:O	3:D:1354:LYS:HG3	2.18	0.43
3:D:1418:LYS:HG2	3:D:1419:PRO:HD2	2.01	0.43
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.52	0.43
1:A:218:LEU:HG	1:B:222:LEU:HD11	2.01	0.43
2:C:726:ILE:HD11	2:C:757:GLY:HA3	2.00	0.43
2:C:853:LEU:HB2	2:C:858:MET:CE	2.48	0.43
3:D:171:LEU:HD12	3:D:390:PRO:HG2	2.01	0.43
2:C:512:ARG:CZ	2:C:523:ILE:HG21	2.49	0.43
3:D:1148:VAL:HA	3:D:1164:ARG:O	2.19	0.43
5:F:364:ARG:HG3	5:F:390:PHE:CE2	2.54	0.43
1:B:72:LYS:HG2	1:B:131:THR:OG1	2.18	0.42
2:C:51:THR:O	2:C:265:ARG:NH2	2.52	0.42
2:C:595:LEU:HD11	2:C:623:TYR:HB3	2.00	0.42
3:D:1107:VAL:HA	3:D:1200:VAL:O	2.18	0.42
3:D:1225:ALA:O	3:D:1229:ILE:HG13	2.18	0.42
3:D:1340:GLY:O	3:D:1344:VAL:HG23	2.18	0.42
3:D:103:TRP:CZ2	3:D:1444:THR:HG22	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:475:LYS:O	3:D:479:GLU:HG2	2.18	0.42
3:D:654:LYS:O	3:D:658:LEU:HG	2.19	0.42
3:D:167:GLU:OE2	3:D:198:ARG:NH1	2.52	0.42
3:D:367:ILE:HG22	3:D:368:VAL:HG23	2.01	0.42
3:D:483:HIS:CE1	3:D:488:ARG:HD3	2.54	0.42
1:A:97:VAL:HG12	1:A:99:LEU:HD12	2.00	0.42
2:C:177:GLU:HG3	2:C:178:PRO:HD2	2.01	0.42
2:C:420:ARG:HD3	2:C:420:ARG:HA	1.83	0.42
2:C:612:VAL:HG22	2:C:622:GLU:HG3	2.00	0.42
5:F:372:ARG:HG2	5:F:386:VAL:HG21	2.01	0.42
5:F:326:ASP:HB3	6:G:19:DG:H22	1.85	0.42
1:A:133:GLU:HG2	1:A:134:GLU:N	2.34	0.42
1:B:124:ASN:N	1:B:124:ASN:OD1	2.52	0.42
3:D:181:ASP:HB2	3:D:205:TYR:CD1	2.54	0.42
3:D:38:LYS:HD3	3:D:38:LYS:HA	1.79	0.42
3:D:41:ARG:HE	3:D:48:ARG:CZ	2.33	0.42
3:D:530:VAL:HG12	3:D:531:ASP:H	1.84	0.42
5:F:276:ARG:O	5:F:279:GLN:HG3	2.19	0.42
3:D:1379:VAL:HG21	3:D:1400:VAL:HG11	2.00	0.42
2:C:195:LEU:O	2:C:199:VAL:HG23	2.20	0.42
2:C:109:LYS:HE2	2:C:368:THR:HG22	2.02	0.42
1:B:85:LEU:HG	1:B:87:VAL:HG23	2.01	0.42
2:C:239:PHE:CD2	2:C:253:ALA:HA	2.54	0.42
2:C:468:ARG:HA	2:C:486:MET:O	2.18	0.42
3:D:1305:LEU:HD13	3:D:1309:ALA:HB3	2.01	0.42
3:D:968:ASP:OD1	3:D:1058:ARG:NH2	2.47	0.42
2:C:1066:ALA:O	2:C:1070:ILE:HG13	2.20	0.42
2:C:710:ILE:HD12	2:C:790:LEU:HB2	2.00	0.42
3:D:1147:ARG:HD3	3:D:1188:VAL:HG11	2.01	0.42
1:B:185:ARG:NH1	1:B:187:GLY:O	2.53	0.42
2:C:76:PRO:HA	2:C:77:PRO:HD2	1.97	0.42
3:D:433:GLY:HA2	3:D:449:SER:N	2.34	0.42
3:D:236:TYR:CD1	3:D:242:LEU:HD12	2.55	0.41
3:D:959:GLU:HB3	3:D:963:TYR:HE1	1.85	0.41
3:D:1205:TYR:O	3:D:1366:LYS:HD3	2.19	0.41
2:C:127:PHE:O	2:C:133:ASP:HA	2.20	0.41
2:C:223:ASP:OD1	2:C:225:SER:OG	2.36	0.41
3:D:963:TYR:CE2	3:D:1002:LYS:HD3	2.55	0.41
3:D:101:HIS:HB3	3:D:104:PHE:HD2	1.86	0.41
3:D:192:ALA:HB1	3:D:193:PRO:HD2	2.01	0.41
2:C:764:GLU:OE2	3:D:54:LYS:HE2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1492:LEU:HD22	4:E:74:VAL:HG21	2.02	0.41
5:F:81:VAL:O	5:F:85:LEU:HG	2.20	0.41
2:C:237:ARG:O	2:C:241:LEU:HG	2.19	0.41
2:C:390:GLN:HB3	2:C:415:PRO:HD3	2.02	0.41
3:D:397:LYS:O	3:D:448:GLU:HG2	2.21	0.41
3:D:587:ARG:NH2	11:D:2110:HOH:O	2.43	0.41
3:D:818:ARG:HE	3:D:820:GLU:CD	2.23	0.41
2:C:441:VAL:HG23	2:C:559:LEU:HA	2.01	0.41
5:F:120:THR:HG22	5:F:122:LEU:HD13	2.03	0.41
2:C:1090:LYS:HD3	2:C:1090:LYS:HA	1.81	0.41
2:C:557:ARG:HG3	2:C:844:GLY:HA3	2.02	0.41
3:D:299:GLU:O	3:D:302:GLN:HG2	2.20	0.41
3:D:34:TYR:CZ	3:D:35:ARG:HG3	2.56	0.41
3:D:959:GLU:HB3	3:D:963:TYR:CE1	2.56	0.41
5:F:236:SER:OG	8:H:5:DA:OP2	2.35	0.41
2:C:27:ARG:NH2	2:C:27:ARG:HB3	2.36	0.41
1:B:111:ALA:HB3	1:B:125:PRO:HA	2.03	0.41
1:A:218:LEU:HD23	1:B:222:LEU:HD21	2.03	0.41
2:C:912:PRO:O	2:C:916:GLU:HG3	2.20	0.41
3:D:840:LYS:HE3	3:D:841:TYR:CZ	2.56	0.41
2:C:571:LEU:HD22	2:C:700:TYR:HA	2.02	0.41
2:C:6:PHE:CD2	2:C:909:ALA:HB2	2.56	0.41
3:D:319:ALA:HA	3:D:337:LEU:HD23	2.03	0.41
6:G:7:DT:H2"	6:G:8:DC:C5	2.56	0.41
1:A:102:LYS:HB2	1:A:102:LYS:HE3	1.84	0.41
1:A:55:SER:HB3	1:A:143:ARG:HB3	2.03	0.41
1:A:6:LEU:HD21	1:A:27:PRO:HG2	2.03	0.40
1:A:183:ASP:HA	2:C:938:LYS:HE3	2.02	0.40
3:D:134:VAL:CG2	3:D:151:GLN:H	2.30	0.40
1:A:11:PHE:O	1:B:228:PRO:HA	2.21	0.40
1:B:80:LEU:HB3	3:D:867:ARG:HH21	1.86	0.40
1:B:90:LEU:HD12	1:B:119:ASP:HA	2.03	0.40
3:D:202:VAL:HG21	3:D:400:VAL:HG13	2.04	0.40
3:D:236:TYR:HB2	3:D:319:ALA:HB3	2.03	0.40
3:D:273:ARG:HB2	3:D:275:GLU:O	2.22	0.40
2:C:274:ARG:HD2	2:C:288:ARG:HG2	2.03	0.40
3:D:1144:LEU:O	3:D:1147:ARG:HG3	2.21	0.40
3:D:643:GLY:HA3	3:D:727:GLN:HB2	2.02	0.40
3:D:849:ALA:O	3:D:853:VAL:HG23	2.22	0.40
2:C:15:LEU:CD1	2:C:583:LEU:HD21	2.51	0.40
2:C:675:ALA:HB2	2:C:867:VAL:HG11	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:658:LEU:HD23	3:D:661:MET:HE1	2.03	0.40
1:A:99:LEU:HB3	1:A:114:PHE:CD2	2.57	0.40
2:C:397:GLU:HG3	2:C:631:SER:HB2	2.03	0.40
2:C:808:ARG:HG3	11:C:1250:HOH:O	2.20	0.40
3:D:222:GLY:HA2	3:D:333:LEU:O	2.21	0.40
3:D:180:LYS:NZ	3:D:357:GLU:OE1	2.54	0.40
3:D:961:LYS:HB2	3:D:961:LYS:HE3	1.90	0.40
5:F:102:LEU:O	5:F:106:VAL:HG23	2.21	0.40
5:F:109:GLY:O	5:F:113:ILE:HG13	2.21	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%)	224 (98%)	5 (2%)	0	100	100
1	B	225/315 (71%)	222 (99%)	3 (1%)	0	100	100
2	C	1108/1119 (99%)	1087 (98%)	21 (2%)	0	100	100
3	D	1483/1524 (97%)	1452 (98%)	31 (2%)	0	100	100
4	E	93/99 (94%)	90 (97%)	3 (3%)	0	100	100
5	F	344/443 (78%)	340 (99%)	4 (1%)	0	100	100
All	All	3482/3815 (91%)	3415 (98%)	67 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	200/273 (73%)	188 (94%)	12 (6%)	19	49
1	B	200/273 (73%)	195 (98%)	5 (2%)	47	75
2	C	936/941 (100%)	884 (94%)	52 (6%)	21	51
3	D	1250/1279 (98%)	1212 (97%)	38 (3%)	41	70
4	E	84/88 (96%)	82 (98%)	2 (2%)	49	75
5	F	301/388 (78%)	288 (96%)	13 (4%)	29	61
All	All	2971/3242 (92%)	2849 (96%)	122 (4%)	30	63

All (122) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	66	SER
1	A	76	VAL
1	A	94	LEU
1	A	96	THR
1	A	112	ARG
1	A	126	ASP
1	A	137	ARG
1	A	148	VAL
1	A	170	VAL
1	A	186	LEU
1	A	205	VAL
1	B	94	LEU
1	B	96	THR
1	B	113	ASP
1	B	143	ARG
1	B	154	GLU
2	C	1	MET
2	C	11	GLU
2	C	27	ARG
2	C	64	LEU
2	C	81	ASP
2	C	100	LEU
2	C	141	HIS
2	C	149	THR
2	C	177	GLU
2	C	182	VAL

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Mol	Chain	Res	Type
2	C	186	VAL
2	C	200	LEU
2	C	203	ASP
2	C	206	THR
2	C	230	ARG
2	C	242	LEU
2	C	317	VAL
2	C	321	GLU
2	C	322	VAL
2	C	348	LEU
2	C	372	LEU
2	C	384	GLU
2	C	402	SER
2	C	418	LEU
2	C	420	ARG
2	C	427	VAL
2	C	430	VAL
2	C	441	VAL
2	C	445	GLU
2	C	454	SER
2	C	455	LEU
2	C	483	VAL
2	C	524	VAL
2	C	578	VAL
2	C	583	LEU
2	C	595	LEU
2	C	610	ARG
2	C	617	ASP
2	C	630	ARG
2	C	657	ASP
2	C	666	LEU
2	C	670	GLN
2	C	674	VAL
2	C	715	THR
2	C	775	ARG
2	C	869	VAL
2	C	878	SER
2	C	907	ASP
2	C	913	GLU
2	C	1075	ASP
2	C	1078	GLU
2	C	1095	LEU

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Mol	Chain	Res	Type
3	D	67	ARG
3	D	80	VAL
3	D	81	THR
3	D	183	GLU
3	D	185	VAL
3	D	199	LEU
3	D	216	VAL
3	D	273	ARG
3	D	275	GLU
3	D	276	ASP
3	D	331	VAL
3	D	335	LEU
3	D	354	VAL
3	D	431	VAL
3	D	503	LEU
3	D	623	VAL
3	D	709	HIS
3	D	717	GLN
3	D	754	PHE
3	D	778	LEU
3	D	783	ARG
3	D	784	ASP
3	D	829	VAL
3	D	833	GLU
3	D	894	LYS
3	D	899	LEU
3	D	904	VAL
3	D	952	ASP
3	D	1041	LEU
3	D	1062	ARG
3	D	1101	VAL
3	D	1129	THR
3	D	1184	GLN
3	D	1188	VAL
3	D	1200	VAL
3	D	1307	LYS
3	D	1312	LEU
3	D	1468	LEU
4	E	50	THR
4	E	93	TYR
5	F	81	VAL
5	F	88	ILE

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Mol	Chain	Res	Type
5	F	94	LEU
5	F	96	LEU
5	F	120	THR
5	F	136	LEU
5	F	209	PHE
5	F	338	LEU
5	F	349	LEU
5	F	358	LEU
5	F	369	LEU
5	F	375	LEU
5	F	392	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	I	0/2	-	-

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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%)	-0.36	3 (1%) 77 60	60, 79, 110, 170	0
1	B	227/315 (72%)	-0.27	3 (1%) 77 60	62, 92, 124, 153	0
2	C	1112/1119 (99%)	-0.30	7 (0%) 89 79	44, 77, 136, 165	0
3	D	1486/1524 (97%)	-0.20	20 (1%) 77 60	41, 79, 142, 177	0
4	E	95/99 (95%)	-0.39	0 100 100	53, 80, 119, 129	0
5	F	346/443 (78%)	-0.17	4 (1%) 79 63	56, 98, 146, 161	0
6	G	16/16 (100%)	-0.60	0 100 100	68, 102, 180, 185	0
7	I	1/2 (50%)	-0.41	0 100 100	72, 72, 72, 72	1 (100%)
8	H	24/24 (100%)	-0.63	0 100 100	92, 126, 177, 189	0
All	All	3538/3857 (91%)	-0.25	37 (1%) 82 69	41, 83, 141, 189	1 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	LEU	4.3
2	C	365	ASP	4.2
2	C	766	GLU	4.1
1	A	232	ALA	3.9
1	A	234	ALA	3.7
3	D	1294	VAL	3.1
5	F	422	LEU	3.1
1	A	233	VAL	3.0
3	D	173	PRO	3.0
3	D	1499	ARG	3.0
2	C	207	LEU	2.9
3	D	409	VAL	2.9
3	D	821	VAL	2.7
3	D	316	GLN	2.6
5	F	149	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
3	D	212	ARG	2.5
3	D	235	ALA	2.5
3	D	1297	GLU	2.4
3	D	345	TYR	2.4
5	F	145	PRO	2.4
3	D	236	TYR	2.4
3	D	1299	PHE	2.4
1	B	3	ASP	2.3
1	B	5	LYS	2.3
2	C	176	VAL	2.3
3	D	213	VAL	2.2
2	C	221	LEU	2.2
2	C	366	SER	2.2
3	D	191	LEU	2.2
3	D	336	PHE	2.1
3	D	974	ILE	2.1
2	C	729	LEU	2.1
3	D	1127	GLU	2.1
3	D	241	ILE	2.1
3	D	1253	THR	2.1
3	D	1300	SER	2.1
5	F	423	ASP	2.1

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 monosaccharides 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(\AA^2)	Q<0.9
9	MG	B	2002	1/1	0.59	0.42	75,75,75,75	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
9	MG	D	2004	1/1	0.88	0.13	78,78,78,78	0
9	MG	B	2001	1/1	0.93	0.08	100,100,100,100	0
9	MG	F	2001	1/1	0.94	0.12	90,90,90,90	0
9	MG	D	2003	1/1	0.97	0.13	46,46,46,46	0
10	ZN	D	2001	1/1	0.98	0.25	99,99,99,99	0
10	ZN	D	2002	1/1	1.00	0.05	120,120,120,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.