



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 18, 2026 – 01:01 AM UTC

PDB ID : 7MLJ / pdb_00007mlj
Title : Crystal structure of Thermus thermophilus reiterative transcription complex with 4nt oligo-G RNA
Authors : Liu, Y.; Ebright, R.H.
Deposited on : 2021-04-28
Resolution : 3.75 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
Xtrriage (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

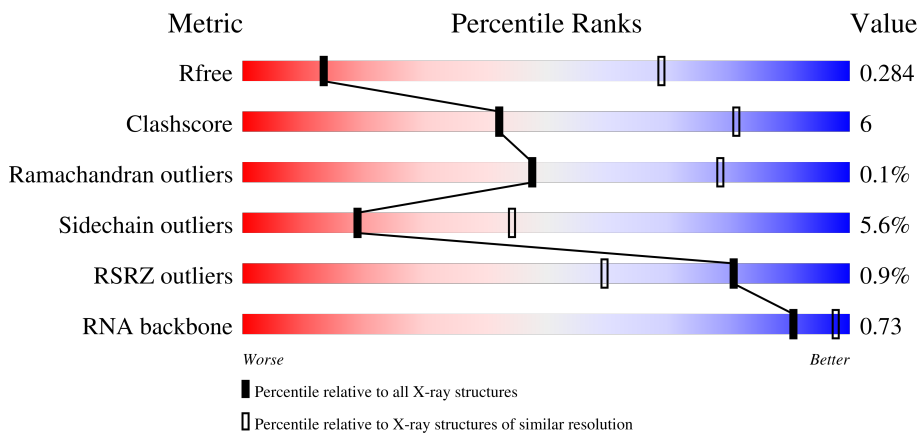
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.75 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	1029 (3.90-3.62)
Clashscore	190562	1061 (3.90-3.62)
Ramachandran outliers	187476	1014 (3.90-3.62)
Sidechain outliers	187428	1009 (3.90-3.62)
RSRZ outliers	180081	1028 (3.90-3.62)
RNA backbone	3983	1001 (4.40-3.10)

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	 58% 14% 28%
1	B	315	 57% 12% 30%
2	C	1119	 80% 18% 2% 2%
3	D	1524	 77% 19% 4% 2%

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Mol	Chain	Length	Quality of chain
4	E	99	 80% 15% 5%
5	F	443	 % 63% 14% 22%
6	G	20	 15% 60% 25%
7	I	4	 50% 50%
8	H	27	 41% 44% 15%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 28489 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
			Total	C	N	O	S			
1	A	226	Total	C	N	O	S	0	0	0
			1782	1138	310	332	2			
1	B	222	Total	C	N	O	S	0	0	0
			1750	1118	304	326	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	1111	Total	C	N	O	S	0	0	0
			8770	5548	1564	1634	24			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	D	1486	Total	C	N	O	S	0	0	0
			11738	7441	2067	2195	35			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
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
			Total	C	N	O	S			
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(P*CP*AP*TP*CP*CP*GP*TP*GP*CP*CP*CP*TP*GP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
6	G	15	304	144	54	91	15	0	0	0

- Molecule 7 is a RNA chain called RNA (5'-R(P*GP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
7	I	4	92	40	20	28	4	0	0	0

- Molecule 8 is a DNA chain called DNA (5'-D(*TP*AP*TP*AP*AP*TP*GP*GP*GP*AP*GP*CP*TP*GP*GP*CP*AP*CP*GP*GP*AP*TP*G)-3').

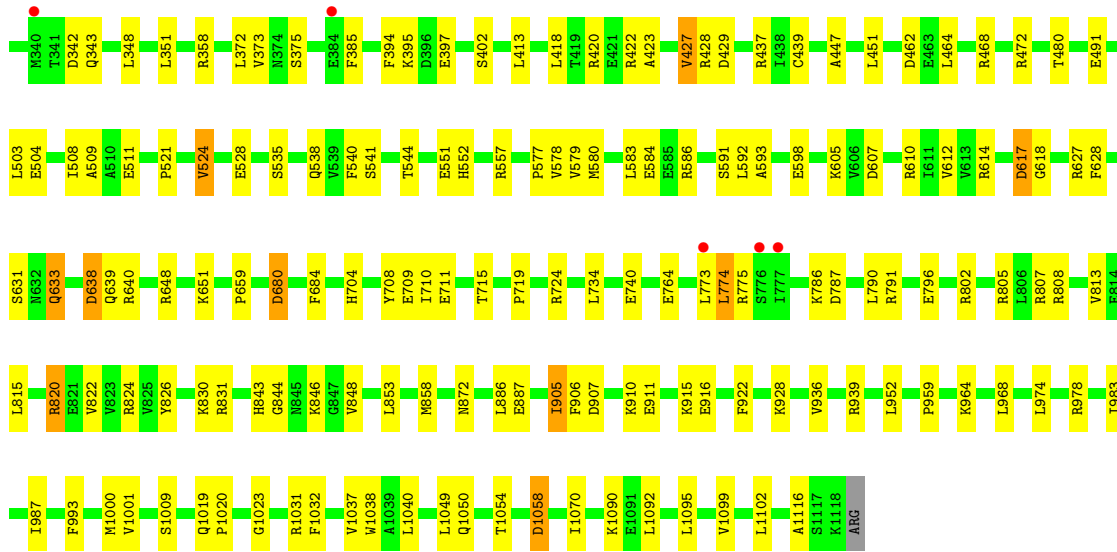
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
8	H	23	478	227	94	135	22	0	0	0

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

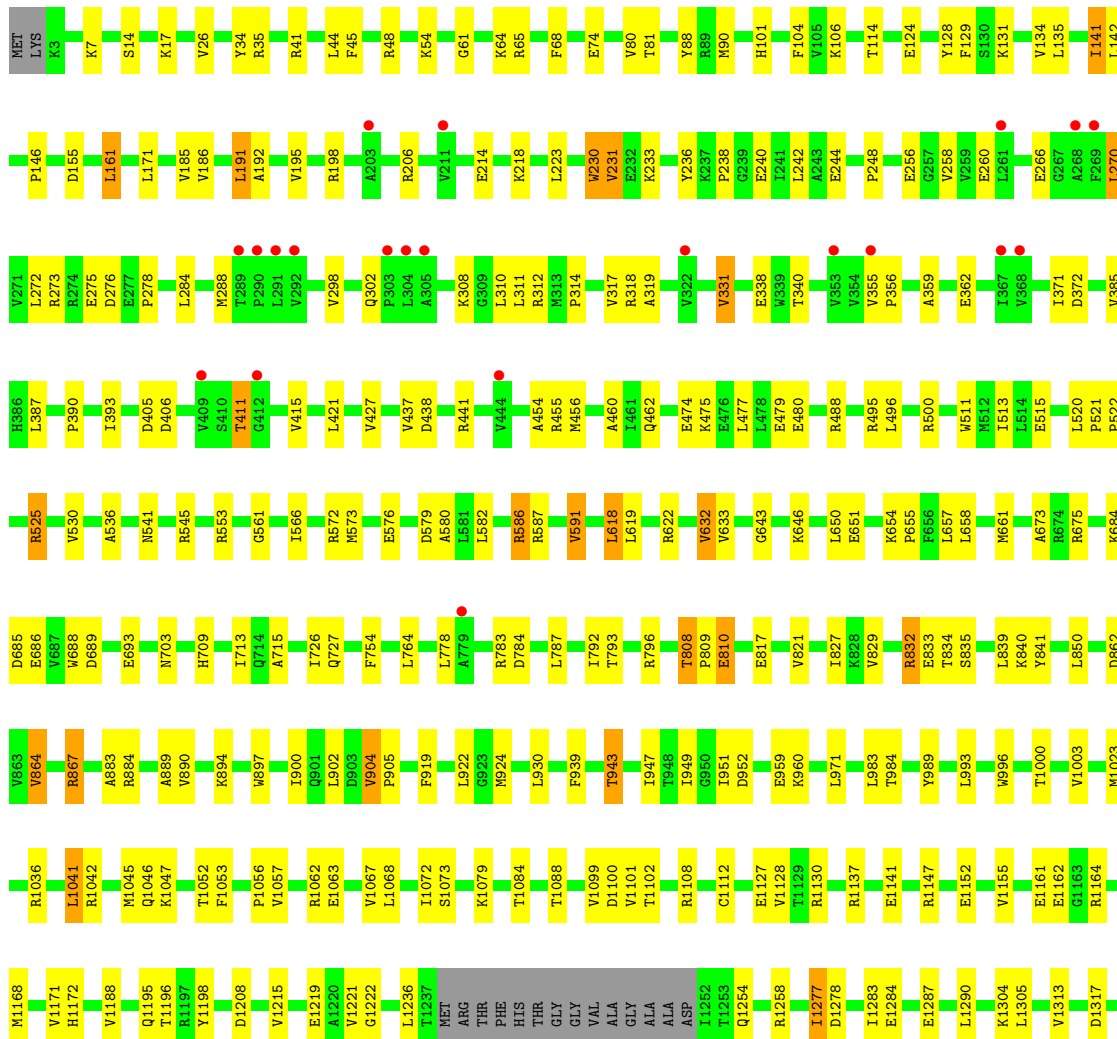
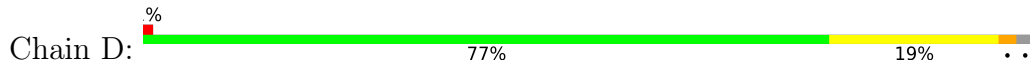
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	2	Total Mg 2 2	0	0
9	D	2	Total Mg 2 2	0	0
9	F	1	Total Mg 1 1	0	0

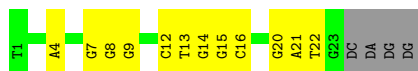
- Molecule 10 is ZINC ION (CCD ID: ZN) (formula: Zn).

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



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





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.31Å 104.66Å 298.41Å 90.00° 98.12° 90.00°	Depositor
Resolution (Å)	46.21 – 3.75 46.21 – 3.75	Depositor EDS
% Data completeness (in resolution range)	97.1 (46.21-3.75) 97.5 (46.21-3.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 3.77Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.265 , 0.283 0.262 , 0.284	Depositor DCC
R_{free} test set	2009 reflections (3.42%)	wwPDB-VP
Wilson B-factor (Å ²)	106.3	Xtrriage
Anisotropy	0.807	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 82.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	28489	wwPDB-VP
Average B, all atoms (Å ²)	152.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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

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	0.11	0/1814	0.31	0/2466
1	B	0.11	0/1782	0.33	0/2424
2	C	0.10	0/8937	0.28	0/12087
3	D	0.10	0/11944	0.28	0/16149
4	E	0.10	0/775	0.25	0/1045
5	F	0.11	0/2852	0.25	0/3837
6	G	0.21	0/339	0.42	0/520
7	I	0.15	0/103	0.26	0/160
8	H	0.22	0/538	0.45	0/831
All	All	0.11	0/29084	0.29	0/39519

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1782	0	1834	22	0
1	B	1750	0	1797	24	0
2	C	8770	0	8874	118	0
3	D	11738	0	11971	165	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	761	0	778	12	0
5	F	2807	0	2882	39	0
6	G	304	0	169	9	0
7	I	92	0	44	2	0
8	H	478	0	260	18	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
All	All	28489	0	28609	359	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 6.

The worst 5 of 359 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)
3:D:1491:THR:HG21	4:E:89:MET:HG2	1.63	0.79
2:C:773:LEU:HB2	5:F:375:LEU:HD11	1.67	0.77
1:B:206:THR:HG22	1:B:209:GLU:H	1.52	0.74
2:C:462:ASP:HB3	2:C:468:ARG:HD2	1.69	0.74
2:C:422:ARG:HA	8:H:16:DC:H5'	1.68	0.74

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	224/315 (71%)	219 (98%)	5 (2%)	0	100 100
1	B	220/315 (70%)	215 (98%)	5 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	1107/1119 (99%)	1082 (98%)	23 (2%)	2 (0%)	43	72
3	D	1482/1524 (97%)	1449 (98%)	31 (2%)	2 (0%)	48	79
4	E	92/99 (93%)	91 (99%)	1 (1%)	0	100	100
5	F	344/443 (78%)	338 (98%)	6 (2%)	0	100	100
All	All	3469/3815 (91%)	3394 (98%)	71 (2%)	4 (0%)	48	79

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	592	LEU
3	D	1440	PHE
2	C	215	GLY
3	D	530	VAL

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	199/273 (73%)	191 (96%)	8 (4%)	28	51
1	B	195/273 (71%)	187 (96%)	8 (4%)	27	51
2	C	936/941 (100%)	879 (94%)	57 (6%)	17	42
3	D	1253/1279 (98%)	1175 (94%)	78 (6%)	16	42
4	E	83/88 (94%)	82 (99%)	1 (1%)	63	71
5	F	301/388 (78%)	286 (95%)	15 (5%)	22	46
All	All	2967/3242 (92%)	2800 (94%)	167 (6%)	19	44

5 of 167 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	D	817	GLU
3	D	1304	LYS
3	D	894	LYS

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Mol	Chain	Res	Type
3	D	1128	VAL
3	D	1501	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
3	D	1172	HIS
3	D	1116	ASN
3	D	350	HIS
3	D	1014	ASN
2	C	860	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
7	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 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	226/315 (71%)	-0.30	0 100 100	113, 141, 165, 179	0
1	B	222/315 (70%)	-0.36	1 (0%) 87 70	105, 148, 179, 195	0
2	C	1111/1119 (99%)	-0.09	6 (0%) 87 70	83, 151, 200, 219	0
3	D	1486/1524 (97%)	-0.10	21 (1%) 73 50	90, 142, 197, 245	1 (0%)
4	E	94/99 (94%)	-0.27	0 100 100	123, 166, 198, 211	0
5	F	346/443 (78%)	0.05	4 (1%) 76 54	112, 160, 213, 223	0
6	G	15/20 (75%)	-0.06	0 100 100	130, 168, 224, 224	0
7	I	4/4 (100%)	0.29	0 100 100	155, 174, 199, 203	0
8	H	23/27 (85%)	0.25	0 100 100	168, 198, 281, 298	0
All	All	3527/3866 (91%)	-0.11	32 (0%) 81 59	83, 149, 201, 298	1 (0%)

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	367	ILE	4.0
3	D	292	VAL	3.8
3	D	290	PRO	3.5
3	D	353	VAL	3.5
3	D	291	LEU	3.4

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(\AA^2)	Q<0.9
9	MG	D	2004	1/1	0.40	0.11	149,149,149,149	0
9	MG	B	2002	1/1	0.75	0.22	92,92,92,92	0
9	MG	F	2001	1/1	0.82	0.09	147,147,147,147	0
9	MG	B	2001	1/1	0.88	0.06	139,139,139,139	0
10	ZN	D	2002	1/1	0.95	0.06	197,197,197,197	0
10	ZN	D	2001	1/1	0.98	0.04	123,123,123,123	0
9	MG	D	2003	1/1	0.98	0.08	98,98,98,98	0

6.5 Other polymers [i](#)

There are no such residues in this entry.