



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 06:29 AM GMT

PDB ID : 1O24
Title : Crystal structure of Thymidylate Synthase Complementing Protein (TM0449) from *Thermotoga maritima* at 2.0 Å resolution
Authors : Mathews, I.I.; Deacon, A.M.; Canaves, J.M.; McMullan, D.; Lesley, S.A.; Agarwalla, S.; Kuhn, P.; Joint Center for Structural Genomics (JCSG)
Deposited on : 2003-02-18
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

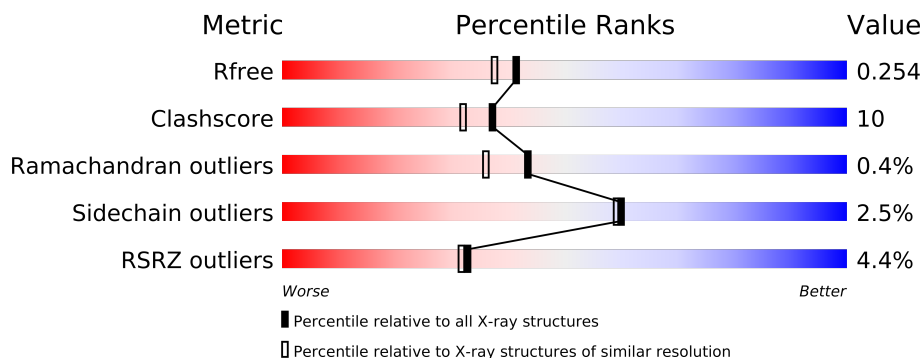
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	232	
1	B	232	
1	C	232	
1	D	232	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7543 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Thymidylate synthase thyX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1743	1136	299	302	6			
1	B	213	Total	C	N	O	S	0	0	0
			1775	1155	306	309	5			
1	C	220	Total	C	N	O	S	0	0	0
			1841	1197	314	324	6			
1	D	214	Total	C	N	O	S	0	0	0
			1776	1156	301	313	6			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	LEADER SEQUENCE	UNP Q9WYT0
A	-10	GLY	-	LEADER SEQUENCE	UNP Q9WYT0
A	-9	SER	-	LEADER SEQUENCE	UNP Q9WYT0
A	-8	ASP	-	LEADER SEQUENCE	UNP Q9WYT0
A	-7	LYS	-	LEADER SEQUENCE	UNP Q9WYT0
A	-6	ILE	-	LEADER SEQUENCE	UNP Q9WYT0
A	-5	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
A	-4	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
A	-3	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
A	-2	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
A	-1	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
A	0	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
B	-11	MET	-	LEADER SEQUENCE	UNP Q9WYT0
B	-10	GLY	-	LEADER SEQUENCE	UNP Q9WYT0
B	-9	SER	-	LEADER SEQUENCE	UNP Q9WYT0
B	-8	ASP	-	LEADER SEQUENCE	UNP Q9WYT0
B	-7	LYS	-	LEADER SEQUENCE	UNP Q9WYT0
B	-6	ILE	-	LEADER SEQUENCE	UNP Q9WYT0
B	-5	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
B	-4	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
B	-3	HIS	-	LEADER SEQUENCE	UNP Q9WYT0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
B	-1	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
B	0	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-11	MET	-	LEADER SEQUENCE	UNP Q9WYT0
C	-10	GLY	-	LEADER SEQUENCE	UNP Q9WYT0
C	-9	SER	-	LEADER SEQUENCE	UNP Q9WYT0
C	-8	ASP	-	LEADER SEQUENCE	UNP Q9WYT0
C	-7	LYS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-6	ILE	-	LEADER SEQUENCE	UNP Q9WYT0
C	-5	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-4	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-3	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-2	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	-1	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
C	0	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-11	MET	-	LEADER SEQUENCE	UNP Q9WYT0
D	-10	GLY	-	LEADER SEQUENCE	UNP Q9WYT0
D	-9	SER	-	LEADER SEQUENCE	UNP Q9WYT0
D	-8	ASP	-	LEADER SEQUENCE	UNP Q9WYT0
D	-7	LYS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-6	ILE	-	LEADER SEQUENCE	UNP Q9WYT0
D	-5	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-4	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-3	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-2	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	-1	HIS	-	LEADER SEQUENCE	UNP Q9WYT0
D	0	HIS	-	LEADER SEQUENCE	UNP Q9WYT0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	96	Total O 96 96	0	0
2	B	107	Total O 107 107	0	0
2	C	115	Total O 115 115	0	0
2	D	90	Total O 90 90	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.48Å 116.64Å 142.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 46.63 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.6 (20.00-2.00) 96.2 (46.63-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.252 0.210 , 0.254	Depositor DCC
R_{free} test set	3016 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	31.1	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 59761 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7543	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.34	0/1791	0.58	0/2422
1	B	0.34	0/1822	0.59	0/2462
1	C	0.35	0/1890	0.59	0/2554
1	D	0.34	0/1824	0.56	0/2467
All	All	0.34	0/7327	0.58	0/9905

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1743	0	1729	37	0
1	B	1775	0	1775	41	0
1	C	1841	0	1839	36	0
1	D	1776	0	1752	41	0
2	A	96	0	0	4	0
2	B	107	0	0	2	0
2	C	115	0	0	4	0
2	D	90	0	0	2	0
All	All	7543	0	7095	144	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (144) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:17:MET:HB2	1:D:17:MET:HB2	1.48	0.93
1:C:106:LEU:HD21	1:C:118:VAL:HG11	1.52	0.92
1:A:17:MET:HB2	1:B:17:MET:HB2	1.52	0.89
1:C:219:GLN:HG3	1:C:220:VAL:HG12	1.64	0.79
1:B:42:ARG:HH11	1:B:42:ARG:HG3	1.49	0.78
1:A:122:ILE:HA	1:A:125:ILE:HD11	1.68	0.74
1:A:163:ASN:ND2	1:A:166:SER:H	1.89	0.71
1:D:192:ARG:HG3	1:D:192:ARG:HH11	1.60	0.67
1:B:163:ASN:ND2	1:B:166:SER:H	1.92	0.67
1:B:106:LEU:CD1	1:B:118:VAL:HB	2.25	0.67
1:A:122:ILE:O	1:A:125:ILE:HD12	1.97	0.65
1:A:125:ILE:HG12	1:D:132:THR:HG21	1.80	0.64
1:A:163:ASN:HD22	1:A:163:ASN:C	2.00	0.64
1:B:42:ARG:HH12	1:B:200:TRP:HA	1.62	0.64
1:A:19:ASN:HB2	2:A:475:HOH:O	1.99	0.63
1:C:192:ARG:HG3	1:C:220:VAL:HG22	1.82	0.62
1:A:167:LEU:HD11	1:A:190:ILE:CG2	2.31	0.61
1:C:195:LYS:HD3	1:C:220:VAL:HG23	1.82	0.61
1:C:125:ILE:HD13	1:C:125:ILE:O	2.01	0.60
1:C:19:ASN:HB2	2:C:477:HOH:O	2.02	0.60
1:A:50:LYS:HG2	1:A:208:TYR:CD2	2.37	0.60
1:D:89:GLY:N	1:D:94:LEU:HD21	2.17	0.59
1:A:46:GLU:O	1:A:50:LYS:HG3	2.02	0.59
1:B:163:ASN:HD22	1:B:163:ASN:C	2.06	0.59
1:B:42:ARG:NH1	1:B:42:ARG:HG3	2.18	0.59
1:B:118:VAL:HG13	1:C:136:LEU:HD22	1.84	0.58
1:D:167:LEU:HD11	1:D:190:ILE:CG2	2.33	0.58
1:A:180:GLN:O	1:A:183:ILE:HG22	2.04	0.58
1:A:167:LEU:HD11	1:A:190:ILE:HG21	1.86	0.57
1:D:180:GLN:O	1:D:183:ILE:HG22	2.04	0.57
1:D:104:GLU:HA	1:D:107:GLU:HG2	1.87	0.57
1:B:127:ASP:O	1:B:131:ARG:HD3	2.04	0.56
1:B:116:GLU:H	1:B:116:GLU:CD	2.09	0.56
1:A:142:PRO:HG3	1:D:109:TYR:CE2	2.40	0.56
1:C:124:GLU:HG2	1:C:128:LYS:HZ2	1.71	0.56
1:D:210:TYR:CE2	1:D:215:LEU:HB2	2.40	0.56
1:A:163:ASN:HD21	1:A:166:SER:H	1.54	0.56
1:A:152:LEU:HD23	1:D:152:LEU:HD23	1.88	0.55
1:C:106:LEU:CD2	1:C:118:VAL:HG11	2.31	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:104:GLU:O	1:A:107:GLU:HB2	2.07	0.55
1:B:192:ARG:HG2	1:B:220:VAL:HG11	1.88	0.55
1:B:192:ARG:HG2	1:B:220:VAL:CG1	2.37	0.55
1:C:121:LYS:HE2	2:C:490:HOH:O	2.06	0.55
1:B:195:LYS:HD3	1:B:220:VAL:HG13	1.88	0.54
1:C:106:LEU:HD22	1:C:118:VAL:HG21	1.89	0.54
1:D:167:LEU:HD11	1:D:190:ILE:HG21	1.90	0.54
1:A:111:THR:HA	1:D:139:SER:O	2.09	0.53
1:B:42:ARG:NH1	2:B:508:HOH:O	2.42	0.52
1:B:106:LEU:HD13	1:B:118:VAL:HB	1.90	0.52
1:D:145:VAL:O	1:D:148:ILE:HG12	2.10	0.52
1:A:87:LEU:HD11	1:A:157:ARG:NH2	2.25	0.52
1:B:106:LEU:HD11	1:B:118:VAL:HB	1.91	0.52
1:A:125:ILE:HD12	1:A:126:VAL:H	1.74	0.52
1:B:111:THR:HA	1:C:139:SER:O	2.10	0.52
1:A:143:ARG:HD3	2:A:341:HOH:O	2.10	0.52
1:D:210:TYR:CZ	1:D:215:LEU:HB2	2.45	0.51
1:C:210:TYR:CE2	1:C:215:LEU:HB2	2.46	0.51
1:C:192:ARG:NH1	1:C:220:VAL:HG13	2.26	0.51
1:A:120:GLU:O	1:A:124:GLU:HG3	2.12	0.50
1:D:29:VAL:O	1:D:29:VAL:HG12	2.12	0.50
1:B:192:ARG:HD3	1:B:192:ARG:C	2.32	0.49
1:C:213:ASP:OD1	1:C:214:ILE:HG23	2.12	0.49
1:D:13:LEU:HD21	1:D:16:VAL:CG2	2.43	0.48
1:B:44:LEU:HD11	1:B:48:LEU:HD11	1.95	0.48
1:C:1:MET:N	2:C:463:HOH:O	2.45	0.48
1:B:192:ARG:HD3	1:B:192:ARG:O	2.13	0.48
1:B:117:ARG:HD2	1:B:121:LYS:HE2	1.95	0.47
1:C:220:VAL:HA	2:C:503:HOH:O	2.13	0.47
1:A:149:VAL:HB	1:D:125:ILE:HG21	1.95	0.47
1:B:42:ARG:NH1	1:B:200:TRP:CD1	2.83	0.47
1:D:54:GLU:HG3	1:D:165:ARG:NE	2.30	0.47
1:A:122:ILE:O	1:A:125:ILE:CD1	2.62	0.47
1:C:106:LEU:HD12	1:C:106:LEU:H	1.80	0.47
1:B:44:LEU:C	1:B:44:LEU:HD13	2.35	0.47
1:C:192:ARG:HH12	1:C:220:VAL:HG13	1.80	0.47
1:D:74:ARG:HD3	2:D:315:HOH:O	2.15	0.47
1:B:106:LEU:O	1:B:107:GLU:C	2.53	0.46
1:B:70:ILE:O	1:B:74:ARG:HG3	2.16	0.46
1:D:126:VAL:HG21	1:D:153:ASN:HD21	1.80	0.46
1:C:106:LEU:O	1:C:109:TYR:HD1	1.99	0.46
1:B:25:ARG:HG3	1:B:30:SER:O	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:55:THR:OG1	1:B:56:PRO:HD3	2.15	0.46
1:C:124:GLU:HG2	1:C:128:LYS:NZ	2.30	0.46
1:D:12:GLU:CG	1:D:65:HIS:HB3	2.46	0.46
1:B:163:ASN:HD21	1:B:166:SER:H	1.62	0.46
1:B:84:TYR:CE2	1:B:160:TRP:CD1	3.03	0.46
1:D:13:LEU:HD11	1:D:16:VAL:CG2	2.45	0.46
1:D:13:LEU:HD21	1:D:16:VAL:HG23	1.98	0.45
1:D:192:ARG:HG3	1:D:192:ARG:NH1	2.28	0.45
1:C:213:ASP:OD1	1:C:214:ILE:N	2.49	0.45
1:C:79:HIS:HE1	1:C:173:LEU:HD12	1.81	0.45
1:D:106:LEU:HD11	1:D:118:VAL:HG11	1.98	0.45
1:D:125:ILE:HD13	1:D:125:ILE:HA	1.86	0.45
1:A:67:LYS:HA	1:A:157:ARG:HG2	1.99	0.45
1:C:104:GLU:O	1:C:107:GLU:HB3	2.16	0.45
1:D:55:THR:N	1:D:56:PRO:CD	2.80	0.45
1:C:192:ARG:HG3	1:C:220:VAL:CG2	2.47	0.45
1:D:103:PRO:HB3	1:D:116:GLU:HG3	1.99	0.45
1:C:88:SER:O	1:C:92:SER:HB2	2.17	0.45
1:D:70:ILE:O	1:D:74:ARG:HG3	2.17	0.44
1:C:33:MET:C	1:C:35:LEU:H	2.20	0.44
1:C:130:TYR:CE2	1:C:134:LEU:HD11	2.52	0.44
1:A:175:ALA:HA	1:A:214:ILE:HD11	1.99	0.44
1:B:192:ARG:HA	1:B:220:VAL:HG12	2.00	0.44
1:C:37:ASP:OD2	1:C:40:ARG:HG2	2.18	0.44
1:C:145:VAL:O	1:C:148:ILE:HG12	2.18	0.44
1:B:202:PHE:CE2	1:B:206:LEU:HD11	2.53	0.43
1:D:62:PHE:O	1:D:161:THR:HA	2.18	0.43
1:A:125:ILE:HD12	1:A:126:VAL:N	2.33	0.43
1:B:122:ILE:O	1:B:126:VAL:HG23	2.19	0.43
1:D:42:ARG:NH1	1:D:200:TRP:CD1	2.79	0.43
1:B:19:ASN:HB2	2:B:504:HOH:O	2.18	0.43
1:D:127:ASP:O	1:D:131:ARG:HD3	2.19	0.43
1:D:144:GLU:HG2	2:D:583:HOH:O	2.18	0.42
1:B:57:PHE:O	1:B:164:ALA:HB3	2.20	0.42
1:A:0:HIS:CE1	2:A:393:HOH:O	2.73	0.42
1:C:133:TYR:O	1:C:137:ILE:HG12	2.20	0.42
1:D:143:ARG:O	1:D:147:ARG:HG2	2.20	0.42
1:A:116:GLU:CD	1:A:116:GLU:H	2.23	0.42
1:A:136:LEU:HD22	1:D:118:VAL:HG13	2.02	0.41
1:D:114:PRO:HA	1:D:115:PRO:HD3	1.97	0.41
1:A:50:LYS:HG2	1:A:208:TYR:CG	2.56	0.41
1:A:50:LYS:NZ	2:A:335:HOH:O	2.53	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:55:THR:OG1	1:C:56:PRO:HD3	2.21	0.41
1:C:62:PHE:O	1:C:161:THR:HA	2.20	0.41
1:A:14:VAL:HB	1:A:63:THR:HG22	2.02	0.41
1:A:114:PRO:HA	1:A:115:PRO:HD3	1.97	0.41
1:B:117:ARG:HH11	1:B:121:LYS:HG3	1.85	0.41
1:A:66:VAL:O	1:A:157:ARG:HA	2.21	0.41
1:D:13:LEU:CD2	1:D:197:LYS:HE3	2.51	0.41
1:A:5:ILE:HD11	1:A:189:ALA:HB2	2.03	0.41
1:D:89:GLY:CA	1:D:94:LEU:HD21	2.51	0.41
1:C:117:ARG:O	1:C:121:LYS:HG2	2.20	0.41
1:D:12:GLU:HG2	1:D:65:HIS:HB3	2.02	0.41
1:B:126:VAL:HG21	1:B:153:ASN:HD21	1.86	0.41
1:B:62:PHE:O	1:B:161:THR:HA	2.21	0.41
1:D:42:ARG:NH1	1:D:200:TRP:CE2	2.89	0.41
1:B:5:ILE:HD11	1:B:189:ALA:HB2	2.03	0.41
1:D:55:THR:OG1	1:D:56:PRO:HD3	2.22	0.40
1:A:0:HIS:HB2	1:A:13:LEU:O	2.21	0.40
1:A:87:LEU:HD12	1:A:157:ARG:HB2	2.04	0.40
1:B:42:ARG:NH1	1:B:200:TRP:CG	2.90	0.40
1:B:142:PRO:HG2	1:C:109:TYR:CE2	2.56	0.40
1:B:1:MET:SD	1:B:196:GLU:OE1	2.79	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	202/232 (87%)	194 (96%)	7 (4%)	1 (0%)	38	29
1	B	209/232 (90%)	203 (97%)	6 (3%)	0	100	100
1	C	218/232 (94%)	208 (95%)	9 (4%)	1 (0%)	38	29
1	D	210/232 (90%)	202 (96%)	7 (3%)	1 (0%)	38	29
All	All	839/928 (90%)	807 (96%)	29 (4%)	3 (0%)	43	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	D	33	MET
1	C	34	GLY

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	185/207 (89%)	178 (96%)	7 (4%)	44	39
1	B	188/207 (91%)	183 (97%)	5 (3%)	57	56
1	C	196/207 (95%)	194 (99%)	2 (1%)	85	88
1	D	187/207 (90%)	182 (97%)	5 (3%)	57	56
All	All	756/828 (91%)	737 (98%)	19 (2%)	60	59

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-1	HIS
1	A	44	LEU
1	A	87	LEU
1	A	105	ARG
1	A	125	ILE
1	A	163	ASN
1	A	192	ARG
1	B	107	GLU
1	B	117	ARG
1	B	127	ASP
1	B	163	ASN
1	B	219	GLN
1	C	104	GLU
1	C	125	ILE
1	D	44	LEU
1	D	87	LEU
1	D	105	ARG
1	D	117	ARG

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Mol	Chain	Res	Type
1	D	125	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	ASN
1	B	163	ASN
1	B	169	ASN
1	B	219	GLN
1	D	75	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	208/232 (89%)	0.18	9 (4%) 34 32	20, 33, 55, 68	0
1	B	213/232 (91%)	0.09	7 (3%) 44 44	21, 33, 62, 72	0
1	C	220/232 (94%)	0.10	12 (5%) 24 23	20, 31, 60, 67	0
1	D	214/232 (92%)	0.24	10 (4%) 30 29	21, 35, 62, 76	0
All	All	855/928 (92%)	0.15	38 (4%) 33 32	20, 33, 61, 76	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	220	VAL	4.6
1	D	109	TYR	4.1
1	D	40	ARG	3.9
1	C	91	TYR	3.8
1	C	138	GLU	3.8
1	D	33	MET	3.7
1	A	32	ASP	3.5
1	A	40	ARG	3.5
1	D	32	ASP	3.3
1	D	34	GLY	3.3
1	D	35	LEU	3.3
1	C	34	GLY	3.1
1	D	106	LEU	3.0
1	A	110	LYS	2.8
1	C	109	TYR	2.8
1	B	112	THR	2.8
1	C	218	VAL	2.7
1	A	33	MET	2.7
1	C	107	GLU	2.7
1	C	33	MET	2.7
1	B	110	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	110	LYS	2.6
1	B	218	VAL	2.4
1	A	107	GLU	2.3
1	B	220	VAL	2.3
1	B	107	GLU	2.3
1	C	94	LEU	2.2
1	D	91	TYR	2.2
1	B	219	GLN	2.2
1	A	-1	HIS	2.1
1	C	40	ARG	2.1
1	A	106	LEU	2.1
1	B	106	LEU	2.1
1	D	107	GLU	2.1
1	A	31	PHE	2.0
1	A	109	TYR	2.0
1	C	35	LEU	2.0
1	C	219	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.