



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:34 AM GMT

PDB ID : 2AVT
Title : Crystal structure of the beta subunit from DNA polymerase of *Streptococcus pyogenes*
Authors : Argiriadi, M.A.; Goedken, E.R.; Bruck, I.; O'donnell, M.; Kuriyan, J.
Deposited on : 2005-08-30
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

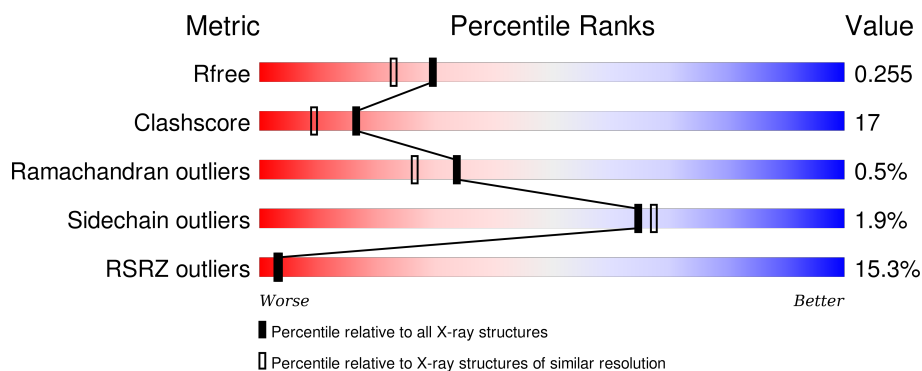
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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. 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	378	
1	B	378	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5943 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 polymerase III beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	0	0
			2899	1845	481	568	5			
1	B	371	Total	C	N	O	S	0	0	0
			2899	1844	483	568	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	PHE	LEU	SEE REMARK 999	GB 16885215
B	199	PHE	LEU	SEE REMARK 999	GB 16885215

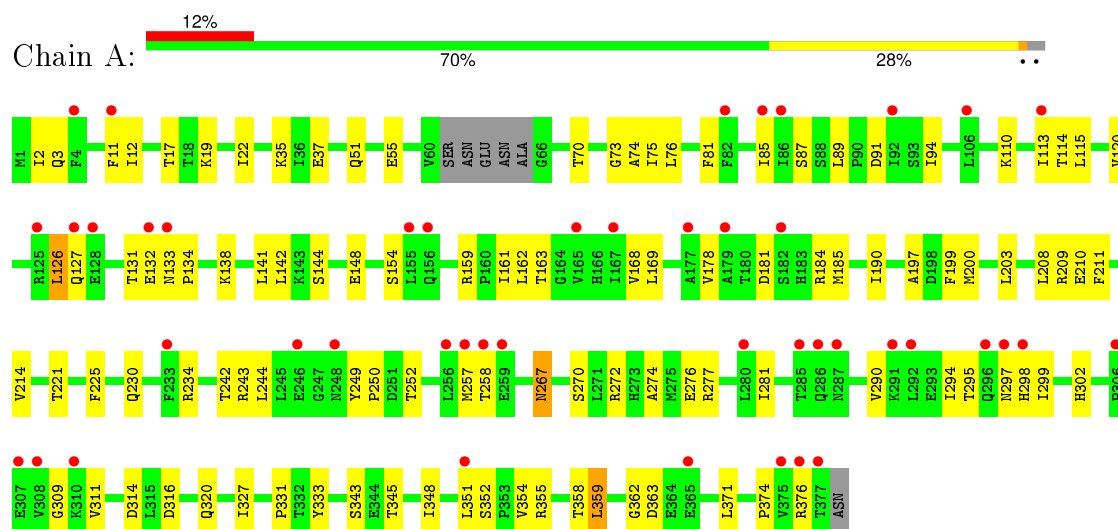
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	65	Total	O	0	0
			65	65		
2	B	80	Total	O	0	0
			80	80		

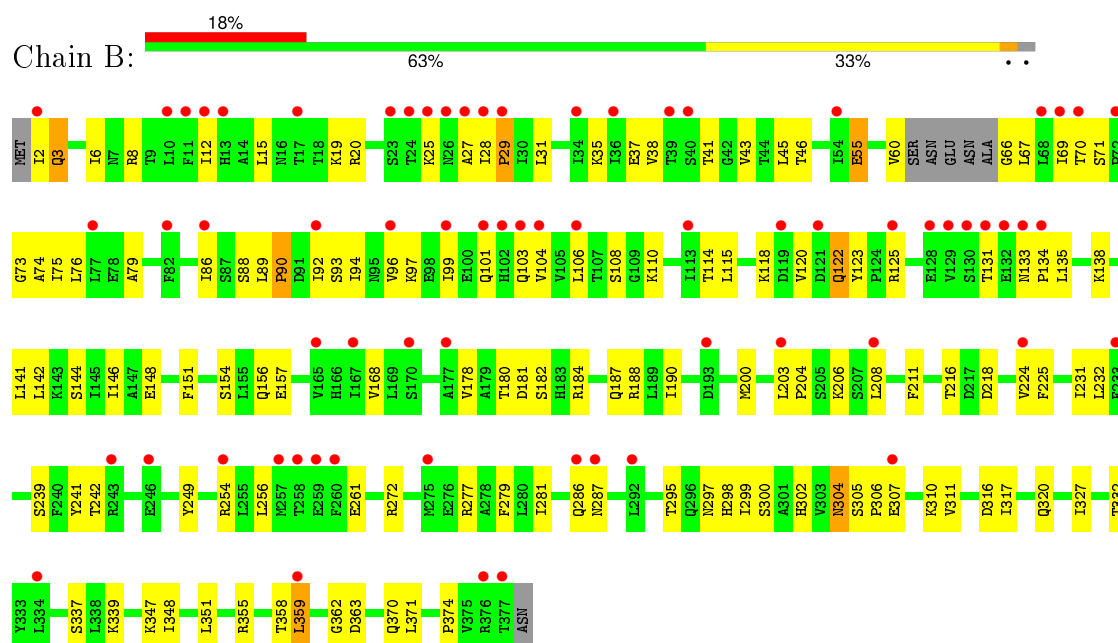
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 polymerase III beta subunit



• Molecule 1: DNA polymerase III beta subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.04Å 74.67Å 82.78Å 90.00° 118.57° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 19.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 90.7 (19.92-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.36 (at 2.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.245 , 0.284 0.249 , 0.255	Depositor DCC
R_{free} test set	2585 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 61.3	EDS
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 52046 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5943	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.33	1/2947 (0.0%)	0.63	0/4003
1	B	0.38	2/2947 (0.1%)	0.64	1/4003 (0.0%)
All	All	0.36	3/5894 (0.1%)	0.64	1/8006 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	157	GLU	CD-OE2	7.32	1.33	1.25
1	A	55	GLU	CD-OE2	7.31	1.33	1.25
1	B	55	GLU	CD-OE2	6.91	1.33	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	359	LEU	CA-CB-CG	5.52	128.00	115.30

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	2899	0	2953	98	0
1	B	2899	0	2957	108	0
2	A	65	0	0	0	0
2	B	80	0	0	2	0
All	All	5943	0	5910	194	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 17.

All (194) 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)
1:B:203:LEU:HD12	1:B:204:PRO:HD2	1.32	1.07
1:A:169:LEU:HB2	1:A:199:PHE:CE2	1.96	0.99
1:B:358:THR:HG22	1:B:371:LEU:HD22	1.45	0.96
1:B:297:ASN:H	1:B:320:GLN:HE22	1.12	0.95
1:A:51:GLN:HE22	1:A:243:ARG:HH21	1.19	0.91
1:B:3:GLN:HE21	1:B:97:LYS:HB2	1.41	0.85
1:A:51:GLN:HE22	1:A:243:ARG:NH2	1.75	0.84
1:B:203:LEU:HD12	1:B:204:PRO:CD	2.06	0.83
1:A:161:ILE:HD12	1:A:162:LEU:N	1.93	0.83
1:B:297:ASN:H	1:B:320:GLN:NE2	1.77	0.81
1:A:3:GLN:HG2	1:A:70:THR:HB	1.64	0.80
1:B:306:PRO:O	1:B:307:GLU:HG2	1.84	0.78
1:A:168:VAL:HG22	1:A:200:MET:HG3	1.66	0.77
1:B:168:VAL:HG22	1:B:200:MET:HG3	1.68	0.76
1:A:203:LEU:HD23	1:A:208:LEU:HD21	1.68	0.75
1:A:297:ASN:HA	1:A:320:GLN:HE22	1.53	0.73
1:A:127:GLN:O	1:A:230:GLN:NE2	2.23	0.71
1:A:169:LEU:CB	1:A:199:PHE:CE2	2.74	0.71
1:A:277:ARG:HE	1:B:90:PRO:HD3	1.55	0.71
1:A:316:ASP:HB2	1:B:110:LYS:NZ	2.07	0.70
1:B:348:ILE:HG12	1:B:359:LEU:HD22	1.73	0.70
1:A:354:VAL:O	1:A:354:VAL:HG12	1.92	0.69
1:A:274:ALA:HA	1:A:277:ARG:NH1	2.09	0.68
1:B:203:LEU:HD23	1:B:208:LEU:HD21	1.76	0.68
1:A:51:GLN:NE2	1:A:243:ARG:HH21	1.90	0.68
1:B:74:ALA:HB1	1:B:120:VAL:HG11	1.76	0.67
1:B:141:LEU:HD11	1:B:190:ILE:CD1	2.25	0.67
1:A:230:GLN:HA	1:A:244:LEU:HG	1.76	0.67
1:A:159:ARG:NE	1:A:161:ILE:HD11	2.10	0.66
1:B:300:SER:OG	1:B:302:HIS:HE1	1.79	0.65
1:A:159:ARG:HD3	1:A:162:LEU:HG	1.79	0.64
1:A:277:ARG:NH2	1:B:90:PRO:HG3	2.12	0.64
1:B:12:ILE:HD11	1:B:86:ILE:HG22	1.79	0.64
1:B:3:GLN:HB3	1:B:70:THR:HB	1.79	0.64
1:B:15:LEU:HD23	1:B:45:LEU:HD13	1.80	0.63
1:A:131:THR:HG22	1:A:131:THR:O	2.00	0.62
1:B:337:SER:HA	1:B:370:GLN:HE22	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:VAL:HG11	1:A:252:THR:HG21	1.81	0.61
1:A:85:ILE:HD13	1:A:115:LEU:HD21	1.82	0.61
1:B:358:THR:CG2	1:B:371:LEU:HD22	2.26	0.61
1:B:90:PRO:HD2	1:B:108:SER:OG	2.01	0.60
1:B:3:GLN:NE2	1:B:97:LYS:HB2	2.14	0.60
1:B:25:LYS:NZ	1:B:25:LYS:HB2	2.17	0.60
1:A:359:LEU:HD22	1:A:359:LEU:N	2.17	0.60
1:A:203:LEU:HD23	1:A:208:LEU:CD2	2.32	0.59
1:A:290:VAL:HB	1:A:331:PRO:HG3	1.84	0.59
1:A:138:LYS:HA	1:A:221:THR:HG22	1.85	0.59
1:B:142:LEU:O	1:B:146:ILE:HG13	2.03	0.58
1:A:297:ASN:HA	1:A:320:GLN:NE2	2.17	0.58
1:A:274:ALA:HA	1:A:277:ARG:HH11	1.67	0.58
1:A:74:ALA:HB1	1:A:120:VAL:HG11	1.85	0.58
1:B:8:ARG:HG3	1:B:94:ILE:HD11	1.86	0.57
1:A:144:SER:O	1:A:148:GLU:HG3	2.04	0.57
1:A:131:THR:HG21	1:A:234:ARG:HH12	1.70	0.57
1:B:144:SER:O	1:B:148:GLU:HG3	2.05	0.57
1:A:154:SER:O	1:A:163:THR:HA	2.04	0.57
1:A:142:LEU:HD23	1:A:211:PHE:CZ	2.40	0.56
1:A:316:ASP:HB2	1:B:110:LYS:HZ3	1.70	0.56
1:A:345:THR:OG1	1:A:362:GLY:HA3	2.06	0.56
1:B:27:ALA:O	1:B:29:PRO:HD3	2.05	0.56
1:B:370:GLN:C	1:B:371:LEU:HD23	2.27	0.55
1:B:131:THR:HG22	1:B:131:THR:O	2.07	0.55
1:B:115:LEU:HD12	1:B:115:LEU:N	2.22	0.55
1:B:231:ILE:HG22	1:B:242:THR:HG23	1.89	0.54
1:B:297:ASN:N	1:B:320:GLN:HE22	1.94	0.54
1:A:272:ARG:O	1:A:276:GLU:HG3	2.08	0.54
1:A:210:GLU:O	1:A:214:VAL:HG23	2.08	0.53
1:A:348:ILE:HG23	1:A:359:LEU:CD1	2.38	0.53
1:B:134:PRO:HB3	1:B:225:PHE:CE2	2.42	0.53
1:A:348:ILE:HG12	1:A:359:LEU:HD12	1.89	0.53
1:A:352:SER:OG	1:A:355:ARG:HG2	2.09	0.53
1:A:169:LEU:HB3	1:A:199:PHE:CD2	2.44	0.52
1:B:304:ASN:C	1:B:304:ASN:HD22	2.11	0.52
1:B:76:LEU:HD11	1:B:123:TYR:HB2	1.90	0.52
1:A:161:ILE:HD12	1:A:162:LEU:H	1.73	0.52
1:A:277:ARG:CZ	1:B:90:PRO:HG3	2.39	0.52
1:A:277:ARG:NE	1:B:90:PRO:HD3	2.23	0.52
1:B:75:ILE:HG12	1:B:76:LEU:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:THR:HB	1:B:310:LYS:HG2	1.91	0.52
1:B:125:ARG:NH1	1:B:125:ARG:HB3	2.25	0.52
1:B:35:LYS:HE2	1:B:37:GLU:CD	2.30	0.52
1:B:231:ILE:CG2	1:B:242:THR:HG23	2.40	0.51
1:B:55:GLU:HB3	1:B:239:SER:HB2	1.91	0.51
1:B:256:LEU:HD13	2:B:426:HOH:O	2.10	0.51
1:A:132:GLU:HG2	1:A:133:ASN:ND2	2.25	0.51
1:A:358:THR:HG22	1:A:371:LEU:CD1	2.41	0.51
1:B:89:LEU:HD12	1:B:94:ILE:HD13	1.93	0.51
1:A:89:LEU:HD12	1:A:94:ILE:HD13	1.92	0.51
1:B:96:VAL:HG22	1:B:106:LEU:CD2	2.41	0.51
1:B:295:THR:OG1	1:B:298:HIS:HB2	2.12	0.50
1:A:203:LEU:HD11	1:A:242:THR:CG2	2.42	0.50
1:B:142:LEU:HD23	1:B:211:PHE:CZ	2.47	0.50
1:A:126:LEU:HD13	1:A:127:GLN:N	2.26	0.50
1:B:35:LYS:HB3	1:B:46:THR:HB	1.94	0.50
1:B:188:ARG:NH2	1:B:190:ILE:HD11	2.27	0.50
1:A:169:LEU:CB	1:A:199:PHE:CD2	2.95	0.49
1:A:181:ASP:OD1	1:A:184:ARG:HG2	2.12	0.49
1:A:294:ILE:HG23	1:A:299:ILE:CD1	2.42	0.49
1:B:3:GLN:HB2	1:B:71:SER:OG	2.11	0.49
1:B:3:GLN:HG2	1:B:97:LYS:HG3	1.93	0.49
1:A:354:VAL:HG22	1:A:376:ARG:HG2	1.94	0.49
1:B:8:ARG:NH2	1:B:90:PRO:O	2.45	0.49
1:A:302:HIS:HA	1:A:311:VAL:O	2.13	0.49
1:A:35:LYS:HE2	1:A:37:GLU:CD	2.32	0.49
1:B:279:PHE:CE1	1:B:332:THR:OG1	2.66	0.49
1:A:169:LEU:HD23	1:A:197:ALA:O	2.13	0.49
1:A:89:LEU:HD21	1:A:113:ILE:HD12	1.95	0.49
1:A:358:THR:HG22	1:A:371:LEU:HD12	1.94	0.48
1:A:184:ARG:HH11	1:A:184:ARG:HG2	1.78	0.48
1:A:134:PRO:HB3	1:A:225:PHE:CE1	2.48	0.48
1:B:38:VAL:HG23	1:B:69:ILE:HD13	1.95	0.48
1:A:127:GLN:HB3	1:A:230:GLN:NE2	2.28	0.48
1:A:309:GLY:HA3	1:B:114:THR:O	2.14	0.48
1:B:43:VAL:HG23	1:B:67:LEU:HD21	1.96	0.48
1:A:81:PHE:CZ	1:A:85:ILE:HD11	2.49	0.48
1:A:348:ILE:HG23	1:A:359:LEU:HD13	1.96	0.48
1:A:184:ARG:CG	1:A:184:ARG:HH11	2.26	0.48
1:B:2:ILE:CD1	1:B:75:ILE:HG22	2.44	0.47
1:B:261:GLU:HG2	1:B:351:LEU:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:ILE:HG13	1:B:317:ILE:HG21	1.96	0.47
1:B:187:GLN:NE2	2:B:401:HOH:O	2.48	0.47
1:A:12:ILE:HD11	1:A:87:SER:HA	1.96	0.47
1:B:327:ILE:HD11	1:B:374:PRO:HB3	1.97	0.47
1:B:138:LYS:HB2	1:B:138:LYS:HE3	1.63	0.46
1:B:304:ASN:HD22	1:B:305:SER:N	2.14	0.46
1:B:118:LYS:HB3	1:B:122:GLN:OE1	2.16	0.46
1:A:249:TYR:CG	1:A:250:PRO:HD2	2.51	0.45
1:A:354:VAL:O	1:A:354:VAL:CG1	2.63	0.45
1:B:2:ILE:HG12	1:B:73:GLY:HA3	1.99	0.45
1:B:3:GLN:HE21	1:B:97:LYS:CB	2.23	0.45
1:A:159:ARG:CD	1:A:161:ILE:HD11	2.46	0.45
1:A:297:ASN:CA	1:A:320:GLN:HE22	2.27	0.45
1:A:19:LYS:HD2	1:A:22:ILE:HB	1.98	0.45
1:B:181:ASP:O	1:B:182:SER:HB2	2.16	0.45
1:B:154:SER:OG	1:B:156:GLN:HG2	2.17	0.44
1:A:2:ILE:HG13	1:A:73:GLY:HA3	1.99	0.44
1:B:232:LEU:HD13	1:B:241:TYR:CZ	2.52	0.44
1:A:267:ASN:HD22	1:A:270:SER:H	1.66	0.44
1:B:178:VAL:HG22	1:B:187:GLN:HG3	2.00	0.44
1:B:92:ILE:HG13	1:B:93:SER:N	2.32	0.44
1:B:89:LEU:HA	1:B:90:PRO:HD3	1.71	0.44
1:A:3:GLN:HG2	1:A:70:THR:CB	2.41	0.43
1:A:110:LYS:HE3	1:A:110:LYS:HB2	1.86	0.43
1:A:290:VAL:CG1	1:A:331:PRO:HG3	2.48	0.43
1:B:281:ILE:HG13	1:B:311:VAL:HG23	1.99	0.43
1:B:286:GLN:O	1:B:287:ASN:HB2	2.18	0.43
1:B:206:LYS:HB2	1:B:206:LYS:NZ	2.33	0.43
1:B:281:ILE:HG13	1:B:311:VAL:CG2	2.48	0.43
1:B:304:ASN:C	1:B:304:ASN:ND2	2.71	0.43
1:B:20:ARG:HG2	1:B:20:ARG:HH11	1.83	0.43
1:A:327:ILE:HD11	1:A:374:PRO:HB3	2.01	0.43
1:B:99:ILE:CG1	1:B:103:GLN:HB2	2.48	0.43
1:B:19:LYS:HG2	1:B:19:LYS:O	2.18	0.43
1:B:151:PHE:HZ	1:B:370:GLN:NE2	2.17	0.43
1:B:347:LYS:HG3	1:B:362:GLY:HA2	2.00	0.43
1:B:272:ARG:HH11	1:B:339:LYS:HG2	1.84	0.43
1:B:203:LEU:CD1	1:B:204:PRO:HD2	2.22	0.43
1:A:351:LEU:HD23	1:A:351:LEU:HA	1.78	0.43
1:B:89:LEU:HB3	1:B:108:SER:OG	2.19	0.42
1:B:135:LEU:HB3	1:B:224:VAL:HB	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:THR:HG22	1:B:249:TYR:OH	2.19	0.42
1:A:343:SER:HB3	1:A:363:ASP:OD2	2.19	0.42
1:A:257:MET:HG2	1:A:358:THR:HG23	1.99	0.42
1:A:267:ASN:ND2	1:A:270:SER:H	2.17	0.42
1:A:89:LEU:HD23	1:A:89:LEU:HA	1.90	0.42
1:A:295:THR:OG1	1:A:298:HIS:HB3	2.20	0.42
1:B:75:ILE:HD13	1:B:104:VAL:CG2	2.50	0.42
1:A:277:ARG:NE	1:B:88:SER:O	2.52	0.42
1:B:28:ILE:HB	1:B:31:LEU:HD12	2.02	0.42
1:A:113:ILE:HD11	1:B:277:ARG:HG2	2.02	0.41
1:A:294:ILE:HG12	1:A:299:ILE:HD12	2.02	0.41
1:A:75:ILE:HG12	1:A:76:LEU:N	2.35	0.41
1:B:348:ILE:CG1	1:B:359:LEU:HD22	2.47	0.41
1:B:216:THR:OG1	1:B:218:ASP:OD1	2.35	0.41
1:A:126:LEU:HD13	1:A:127:GLN:H	1.86	0.41
1:A:17:THR:CG2	1:A:214:VAL:HG13	2.51	0.41
1:B:2:ILE:HG23	1:B:38:VAL:HG11	2.02	0.41
1:A:281:ILE:HG13	1:A:311:VAL:CG2	2.50	0.41
1:A:131:THR:HG21	1:A:234:ARG:NH1	2.34	0.41
1:A:110:LYS:NZ	1:B:316:ASP:OD2	2.54	0.41
1:B:355:ARG:HG3	1:B:355:ARG:HH11	1.86	0.41
1:B:370:GLN:O	1:B:371:LEU:HD23	2.21	0.41
1:B:211:PHE:CD1	1:B:211:PHE:C	2.94	0.41
1:A:316:ASP:HB2	1:B:110:LYS:HZ2	1.81	0.40
1:A:185:MET:SD	1:A:252:THR:OG1	2.79	0.40
1:A:290:VAL:CB	1:A:331:PRO:HG3	2.50	0.40
1:A:184:ARG:HD3	1:A:333:TYR:CD1	2.56	0.40
1:B:133:ASN:HA	1:B:134:PRO:HD2	1.81	0.40
1:B:35:LYS:HE2	1:B:37:GLU:CG	2.51	0.40
1:B:101:GLN:HB2	1:B:103:GLN:HE21	1.86	0.40
1:B:99:ILE:HG12	1:B:103:GLN:HB2	2.03	0.40
1:B:41:THR:O	1:B:60:VAL:HG13	2.20	0.40
1:B:6:ILE:HG13	1:B:66:GLY:O	2.21	0.40
1:A:141:LEU:HD11	1:A:190:ILE:CD1	2.52	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	368/378 (97%)	349 (95%)	18 (5%)	1 (0%)	46	41
1	B	367/378 (97%)	354 (96%)	10 (3%)	3 (1%)	24	15
All	All	735/756 (97%)	703 (96%)	28 (4%)	4 (0%)	34	26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	THR
1	B	79	ALA
1	B	90	PRO
1	B	29	PRO

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	339/346 (98%)	332 (98%)	7 (2%)	61	63
1	B	340/346 (98%)	334 (98%)	6 (2%)	66	69
All	All	679/692 (98%)	666 (98%)	13 (2%)	65	67

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	PHE
1	A	91	ASP

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Mol	Chain	Res	Type
1	A	126	LEU
1	A	209	ARG
1	A	267	ASN
1	A	314	ASP
1	A	359	LEU
1	B	3	GLN
1	B	122	GLN
1	B	184	ARG
1	B	254	ARG
1	B	304	ASN
1	B	363	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	102	HIS
1	A	133	ASN
1	A	267	ASN
1	A	320	GLN
1	B	3	GLN
1	B	95	ASN
1	B	103	GLN
1	B	166	HIS
1	B	187	GLN
1	B	302	HIS
1	B	304	ASN
1	B	320	GLN
1	B	370	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	372/378 (98%)	0.81	45 (12%) 6 6	18, 40, 62, 75	0
1	B	371/378 (98%)	0.95	69 (18%) 2 2	17, 38, 62, 76	0
All	All	743/756 (98%)	0.88	114 (15%) 3 3	17, 39, 62, 76	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	131	THR	7.2
1	A	286	GLN	6.8
1	B	258	THR	6.2
1	A	258	THR	6.2
1	A	297	ASN	6.0
1	B	130	SER	5.1
1	B	29	PRO	4.7
1	B	39	THR	4.6
1	B	132	GLU	4.6
1	B	27	ALA	4.5
1	A	298	HIS	4.4
1	B	70	THR	4.3
1	A	351	LEU	4.2
1	A	306	PRO	4.1
1	B	68	LEU	4.1
1	A	287	ASN	4.0
1	A	308	VAL	4.0
1	B	24	THR	4.0
1	B	260	PHE	3.9
1	A	132	GLU	3.6
1	B	13	HIS	3.6
1	A	377	THR	3.6
1	A	256	LEU	3.5
1	A	307	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	296	GLN	3.3
1	B	177	ALA	3.3
1	B	92	ILE	3.3
1	A	133	ASN	3.2
1	B	96	VAL	3.2
1	B	54	ILE	3.2
1	B	106	LEU	3.2
1	B	377	THR	3.2
1	B	119	ASP	3.2
1	B	246	GLU	3.1
1	B	101	GLN	3.1
1	B	10	LEU	3.1
1	B	40	SER	3.0
1	B	34	ILE	3.0
1	A	233	PHE	3.0
1	A	285	THR	3.0
1	B	26	ASN	3.0
1	B	69	ILE	2.9
1	B	359	LEU	2.9
1	B	259	GLU	2.9
1	A	127	GLN	2.9
1	B	243	ARG	2.9
1	A	310	LYS	2.9
1	B	170	SER	2.8
1	A	82	PHE	2.7
1	B	287	ASN	2.7
1	A	125	ARG	2.7
1	A	182	SER	2.7
1	B	23	SER	2.7
1	A	246	GLU	2.7
1	B	307	GLU	2.7
1	B	99	ILE	2.7
1	B	133	ASN	2.6
1	B	334	LEU	2.6
1	B	233	PHE	2.6
1	A	292	LEU	2.6
1	B	25	LYS	2.6
1	B	36	ILE	2.6
1	B	86	ILE	2.6
1	A	248	ASN	2.6
1	B	17	THR	2.6
1	A	259	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	257	MET	2.6
1	B	286	GLN	2.5
1	B	203	LEU	2.5
1	A	165	VAL	2.5
1	B	113	ILE	2.5
1	B	77	LEU	2.5
1	B	103	GLN	2.5
1	B	167	ILE	2.5
1	A	128	GLU	2.5
1	A	156	GLN	2.5
1	B	28	ILE	2.4
1	B	224	VAL	2.4
1	B	193	ASP	2.4
1	A	177	ALA	2.4
1	A	155	LEU	2.4
1	B	82	PHE	2.4
1	B	125	ARG	2.3
1	A	167	ILE	2.3
1	A	106	LEU	2.3
1	B	292	LEU	2.3
1	A	92	ILE	2.3
1	B	104	VAL	2.2
1	B	254	ARG	2.2
1	A	86	ILE	2.2
1	B	102	HIS	2.2
1	A	280	LEU	2.2
1	B	12	ILE	2.2
1	A	375	VAL	2.2
1	B	129	VAL	2.2
1	B	72	PRO	2.2
1	B	121	ASP	2.1
1	A	113	ILE	2.1
1	A	291	LYS	2.1
1	B	11	PHE	2.1
1	B	128	GLU	2.1
1	A	179	ALA	2.1
1	A	376	ARG	2.1
1	B	275	MET	2.1
1	B	2	ILE	2.1
1	A	4	PHE	2.1
1	A	11	PHE	2.1
1	A	85	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	134	PRO	2.0
1	B	208	LEU	2.0
1	A	257	MET	2.0
1	B	376	ARG	2.0
1	B	165	VAL	2.0
1	A	365	GLU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.