



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 01:39 pm BST

PDB ID : 3U58  
Title : Crystal Structure of the Tetrahymena telomerase processivity factor Teb1 AB  
Authors : Zeng, Z.; Huang, J.; Yang, Y.; Lei, M.  
Deposited on : 2011-10-11  
Resolution : 2.61 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](mailto: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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

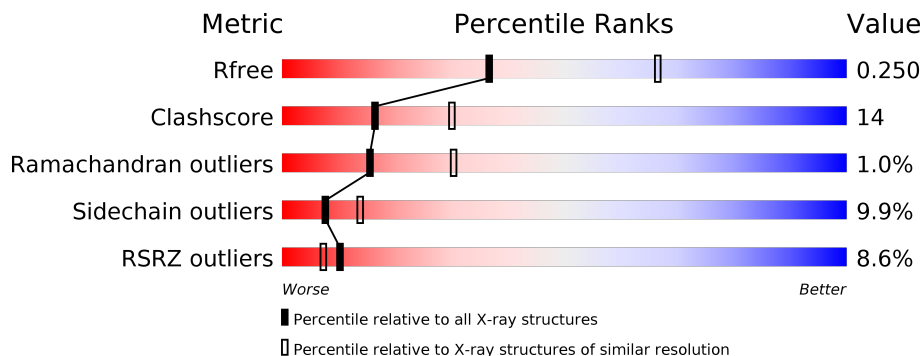
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.61 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.



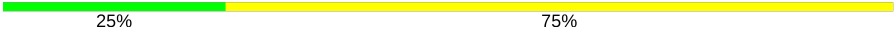

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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  $\geq 3$ , 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	213	<div style="display: flex; align-items: center;"> <div style="width: 10%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 65%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 29%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 7%; height: 10px; background-color: orange;"></div> </div>
1	B	213	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 64%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 29%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: orange;"></div> </div>
1	C	213	<div style="display: flex; align-items: center;"> <div style="width: 9%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 69%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 26%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange;"></div> </div>
1	D	213	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 69%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: orange;"></div> </div>
2	E	4	<div style="display: flex; align-items: center;"> <div style="width: 50%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 50%; height: 10px; background-color: yellow; margin-right: 5px;"></div> </div>
2	F	4	<div style="display: flex; align-items: center;"> <div style="width: 25%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 50%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: orange;"></div> </div>

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Mol	Chain	Length	Quality of chain
2	G	4	 25% 75%
2	H	4	 25% 75%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7378 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 Tetrahymena Teb1 AB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	213	Total 1759	C 1128	N 290	O 338	S 3	0	0	0
1	B	210	Total 1739	C 1118	N 287	O 331	S 3	0	0	0
1	C	211	Total 1743	C 1120	N 288	O 332	S 3	0	0	0
1	D	210	Total 1739	C 1118	N 287	O 331	S 3	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*GP\*GP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	4	Total 83	C 40	N 17	O 23	P 3	0	0	0
2	H	4	Total 83	C 40	N 17	O 23	P 3	0	0	0
2	F	4	Total 83	C 40	N 17	O 23	P 3	0	0	0
2	G	4	Total 83	C 40	N 17	O 23	P 3	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total 15	O 15	0	0
3	B	20	Total 20	O 20	0	0
3	C	11	Total 11	O 11	0	0
3	D	17	Total 17	O 17	0	0

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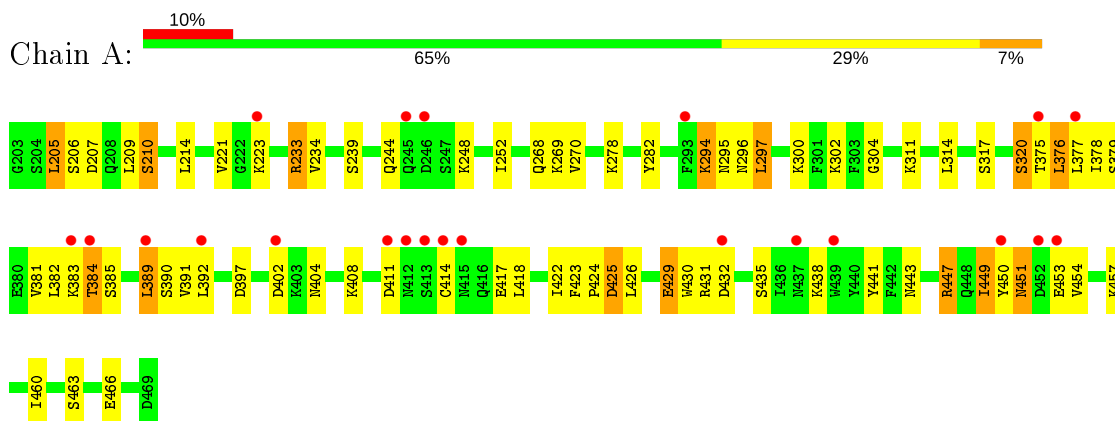
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
3	E	1	Total O 1 1	0	0
3	H	1	Total O 1 1	0	0
3	F	1	Total O 1 1	0	0

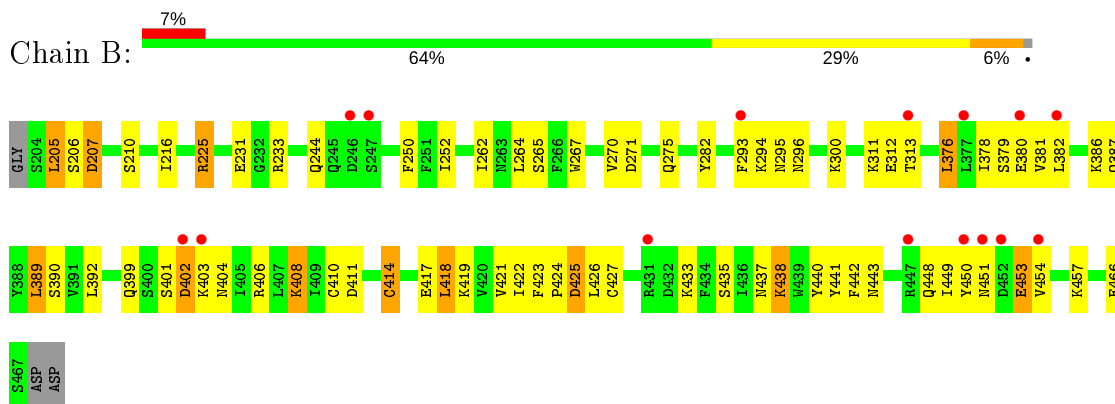
### 3 Residue-property plots

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

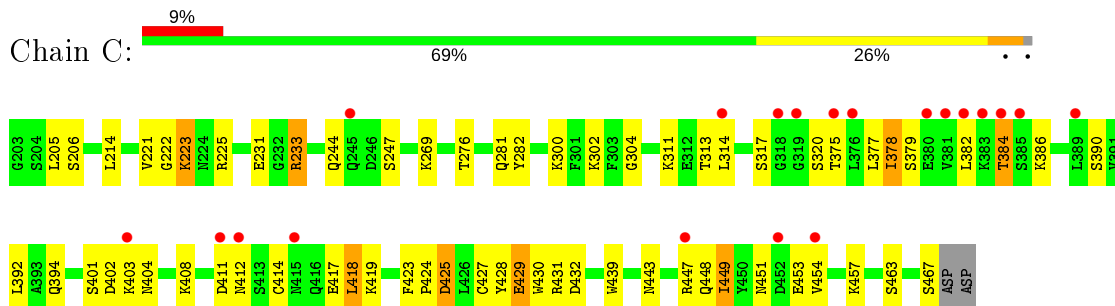
- Molecule 1: Tetrahymena Teb1 AB



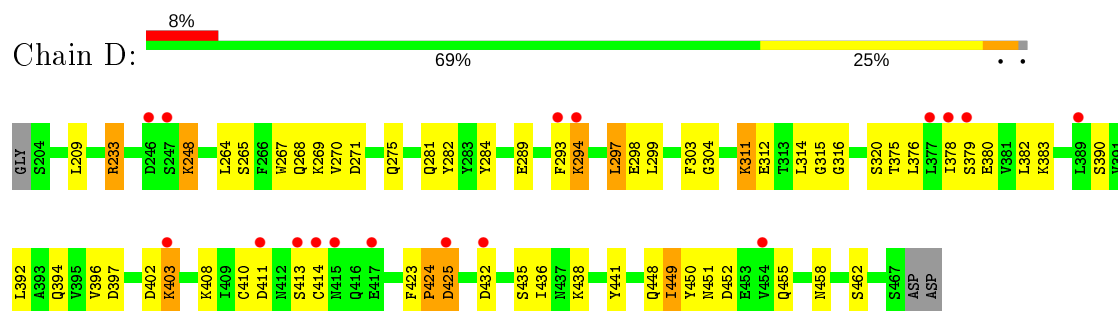
- Molecule 1: Tetrahymena Teb1 AB



- Molecule 1: Tetrahymena Teb1 AB



- Molecule 1: Tetrahymena Teb1 AB



- Molecule 2: DNA (5'-D(\*GP\*GP\*GP\*T)-3')



- Molecule 2: DNA (5'-D(\*GP\*GP\*GP\*T)-3')



- Molecule 2: DNA (5'-D(\*GP\*GP\*GP\*T)-3')



- Molecule 2: DNA (5'-D(\*GP\*GP\*GP\*T)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.05Å 83.11Å 82.89Å 108.45° 111.59° 108.42°	Depositor
Resolution (Å)	33.58 – 2.61 39.34 – 2.61	Depositor EDS
% Data completeness (in resolution range)	94.0 (33.58-2.61) 93.7 (39.34-2.61)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.7_650	Depositor
R, $R_{free}$	0.219 , 0.259 0.211 , 0.250	Depositor DCC
$R_{free}$ test set	4907 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.0	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 27.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage

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<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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.



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Property	Value	Source
Estimated twinning fraction	0.000 for h+k+1,-l,-h	Xtrriage
	0.000 for -l,h+k+1,-k	
	0.429 for h+k+1,-h,-k	
	0.429 for -k,-l,h+k+1	
	0.000 for -k,h+k+1,-h	
	0.000 for -l,-h,h+k+1	
	0.000 for k,-h-k-l,l	
	0.000 for -h-k-l,h,l	
	0.000 for l,h,k	
	0.000 for k,l,h	
	0.000 for h,l,-h-k-l	
	0.000 for h,-h-k-l,k	
	0.000 for l,k,-h-k-l	
	0.000 for -h-k-l,k,h	
	0.019 for k,h,-h-k-l	
	0.000 for -h,-k,h+k+1	
	0.438 for l,-h-k-l,h	
	0.019 for -l,-k,-h	
0.019 for -h-k-l,l,k		
0.000 for h+k+1,-k,-l		
0.000 for -k,-h,-l		
0.018 for -h,h+k+1,-l		
0.000 for -h,-l,-k		
F <sub>o</sub> ,F <sub>c</sub> correlation	0.94	EDS
Total number of atoms	7378	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% of the height of the origin peak. No significant pseudotranslation is detected.*

## 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	0.38	0/1794	0.59	4/2414 (0.2%)
1	B	0.37	0/1774	0.55	2/2387 (0.1%)
1	C	0.37	0/1778	0.54	1/2392 (0.0%)
1	D	0.38	0/1774	0.57	2/2387 (0.1%)
2	E	0.62	0/93	1.00	0/143
2	F	0.70	0/93	1.29	2/143 (1.4%)
2	G	0.60	0/93	1.14	0/143
2	H	0.63	0/93	1.15	0/143
All	All	0.39	0/7492	0.61	11/10152 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	315	GLY	N-CA-C	8.19	133.56	113.10
1	C	404	ASN	N-CA-CB	-6.75	98.44	110.60
2	F	4	DT	N3-C4-O4	6.56	123.84	119.90
1	A	383	LYS	CB-CA-C	-6.22	97.96	110.40
2	F	4	DT	C5-C4-O4	-6.03	120.68	124.90
1	A	384	THR	CB-CA-C	-6.01	95.38	111.60
1	A	385	SER	N-CA-CB	-5.63	102.05	110.50
1	B	437	ASN	CB-CA-C	-5.61	99.18	110.40
1	D	316	GLY	N-CA-C	-5.46	99.45	113.10
1	B	404	ASN	N-CA-CB	-5.27	101.11	110.60
1	A	404	ASN	N-CA-CB	-5.18	101.27	110.60

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	1759	0	1729	54	0
1	B	1739	0	1718	58	0
1	C	1743	0	1721	51	0
1	D	1739	0	1718	43	0
2	E	83	0	47	2	0
2	F	83	0	47	3	0
2	G	83	0	47	2	0
2	H	83	0	47	2	0
3	A	15	0	0	3	0
3	B	20	0	0	4	0
3	C	11	0	0	0	0
3	D	17	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	H	1	0	0	0	0
All	All	7378	0	7074	206	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 14.

All (206) 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:C:378:ILE:HD12	1:C:411:ASP:HB2	1.32	1.10
1:A:378:ILE:HG12	1:A:411:ASP:HB3	1.63	0.81
1:C:403:LYS:O	1:C:424:PRO:HD3	1.82	0.79
1:B:406:ARG:HH11	1:B:419:LYS:HZ3	1.30	0.79
1:D:402:ASP:O	1:D:403:LYS:HB2	1.86	0.75
1:D:423:PHE:HB3	1:D:424:PRO:HD2	1.69	0.75
1:A:376:LEU:H	1:A:376:LEU:HD12	1.51	0.74
1:B:406:ARG:HH11	1:B:419:LYS:NZ	1.85	0.74
1:C:269:LYS:HE2	1:C:304:GLY:HA2	1.71	0.73
2:F:4:DT:H2'	2:F:4:DT:OP2	1.90	0.72
1:A:384:THR:HG23	1:A:384:THR:O	1.90	0.72
1:A:397:ASP:HB3	1:A:408:LYS:HB2	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:424:PRO:O	1:D:425:ASP:HB2	1.91	0.71
1:C:233:ARG:HD3	1:C:282:TYR:CZ	2.25	0.71
1:B:233:ARG:HD3	1:B:282:TYR:CZ	2.26	0.71
1:A:423:PHE:HB3	1:A:424:PRO:HD2	1.72	0.70
1:C:424:PRO:O	1:C:425:ASP:HB2	1.93	0.69
1:C:378:ILE:CD1	1:C:411:ASP:HB2	2.16	0.69
1:C:378:ILE:HD13	1:C:379:SER:H	1.56	0.68
1:B:231:GLU:CD	1:B:311:LYS:HE2	2.15	0.68
1:A:389:LEU:HD13	3:A:3:HOH:O	1.94	0.67
1:D:312:GLU:OE1	2:G:1:DG:H5'	1.94	0.67
1:A:378:ILE:CG1	1:A:411:ASP:HB3	2.25	0.67
1:C:423:PHE:HB3	1:C:424:PRO:HD2	1.76	0.66
1:A:378:ILE:HG13	1:A:379:SER:N	2.09	0.66
1:B:423:PHE:HB3	1:B:424:PRO:HD2	1.78	0.66
1:B:411:ASP:OD2	1:B:414:CYS:HB2	1.96	0.65
1:A:207:ASP:O	1:A:210:SER:HB2	1.96	0.65
1:A:376:LEU:HD13	1:A:381:VAL:HG22	1.78	0.65
1:C:222:GLY:HA2	1:D:233:ARG:HH22	1.61	0.65
1:B:225:ARG:HG3	1:C:282:TYR:OH	1.97	0.64
1:A:378:ILE:HG13	1:A:379:SER:H	1.61	0.64
1:B:424:PRO:O	1:B:425:ASP:HB2	1.96	0.63
1:D:233:ARG:HD3	1:D:282:TYR:CZ	2.33	0.63
1:A:317:SER:O	1:A:320:SER:HB2	1.99	0.62
1:B:389:LEU:HD13	3:B:2:HOH:O	2.00	0.62
1:B:406:ARG:NH1	1:B:419:LYS:NZ	2.48	0.61
1:B:271:ASP:O	1:B:275:GLN:HG2	2.01	0.60
1:D:397:ASP:HB3	1:D:408:LYS:HB2	1.82	0.60
1:A:423:PHE:HE1	1:A:457:LYS:HE2	1.65	0.60
1:B:389:LEU:HB2	3:B:2:HOH:O	2.01	0.59
1:D:435:SER:HB2	1:D:438:LYS:CD	2.32	0.59
2:H:4:DT:H5'	2:H:4:DT:H6	1.66	0.59
1:A:423:PHE:CE1	1:A:457:LYS:HE2	2.38	0.58
1:A:378:ILE:CD1	1:A:411:ASP:HB3	2.34	0.58
1:C:429:GLU:HB3	1:C:430:TRP:CD1	2.39	0.58
1:C:244:GLN:HB2	1:C:247:SER:HB3	1.85	0.58
1:D:448:GLN:HG2	1:D:449:ILE:H	1.68	0.58
1:B:378:ILE:HB	1:B:411:ASP:OD2	2.03	0.57
1:B:435:SER:HB2	1:B:438:LYS:HD3	1.86	0.57
1:A:233:ARG:HD3	1:A:282:TYR:CZ	2.40	0.57
1:C:317:SER:HB2	1:C:320:SER:OG	2.05	0.57
1:C:448:GLN:OE1	1:C:457:LYS:HD3	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:378:ILE:HD12	1:C:411:ASP:CB	2.21	0.56
1:C:424:PRO:HA	1:C:427:CYS:SG	2.45	0.56
1:A:449:ILE:HG12	1:A:454:VAL:HG12	1.88	0.56
1:B:231:GLU:OE2	1:B:311:LYS:HE2	2.06	0.56
1:B:225:ARG:HG3	1:C:282:TYR:CZ	2.40	0.56
1:C:379:SER:HB3	1:C:414:CYS:HB2	1.88	0.55
1:C:423:PHE:CB	1:C:424:PRO:HD2	2.36	0.55
1:A:381:VAL:HG21	1:A:391:VAL:HG22	1.87	0.55
2:F:2:DG:H2''	2:F:3:DG:OP1	2.07	0.55
1:D:396:VAL:O	1:D:436:ILE:HD11	2.07	0.54
1:B:378:ILE:HG13	1:B:411:ASP:HB3	1.90	0.54
1:B:424:PRO:HA	1:B:427:CYS:SG	2.48	0.54
1:C:402:ASP:HA	1:C:431:ARG:NH1	2.22	0.54
1:B:392:LEU:HD13	1:B:441:TYR:CD1	2.43	0.54
1:D:423:PHE:CB	1:D:424:PRO:HD2	2.38	0.54
1:A:384:THR:O	1:A:384:THR:CG2	2.56	0.54
2:H:2:DG:H2''	2:H:3:DG:OP1	2.08	0.53
1:C:231:GLU:CD	1:C:311:LYS:HZ2	2.11	0.53
1:A:392:LEU:HD13	1:A:441:TYR:CE1	2.44	0.53
1:D:396:VAL:C	1:D:436:ILE:HD11	2.29	0.53
1:D:269:LYS:HE2	1:D:304:GLY:HA2	1.91	0.53
1:B:408:LYS:HA	1:B:418:LEU:O	2.09	0.52
1:A:376:LEU:HD11	1:A:389:LEU:HB2	1.91	0.52
1:C:394:GLN:HB3	1:C:439:TRP:CE2	2.45	0.52
1:A:389:LEU:HB2	3:A:3:HOH:O	2.11	0.51
1:B:231:GLU:OE1	1:B:311:LYS:HE2	2.11	0.51
1:D:284:TYR:CZ	1:D:311:LYS:HE2	2.45	0.51
1:A:376:LEU:HD13	1:A:381:VAL:CG2	2.41	0.51
1:B:376:LEU:CD1	1:B:381:VAL:HG22	2.41	0.50
1:B:433:LYS:HG2	1:B:433:LYS:O	2.11	0.50
1:C:378:ILE:HG12	1:C:414:CYS:HB3	1.94	0.50
1:B:392:LEU:HD13	1:B:441:TYR:CE1	2.47	0.50
1:B:451:ASN:H	1:B:453:GLU:CD	2.15	0.50
1:C:402:ASP:HA	1:C:431:ARG:HH12	1.76	0.50
2:E:2:DG:H2''	2:E:3:DG:OP1	2.11	0.50
1:A:375:THR:HA	1:A:390:SER:O	2.12	0.49
1:B:422:ILE:HD13	1:B:422:ILE:N	2.26	0.49
1:D:265:SER:HB2	1:D:267:TRP:NE1	2.27	0.49
1:D:403:LYS:O	1:D:424:PRO:HD3	2.12	0.49
1:B:421:VAL:C	1:B:422:ILE:HD13	2.33	0.49
1:D:293:PHE:O	1:D:294:LYS:HB2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:GLN:HG3	1:C:247:SER:HB3	1.94	0.49
1:C:375:THR:HA	1:C:390:SER:O	2.12	0.49
1:B:205:LEU:HD23	1:B:206:SER:H	1.77	0.48
1:D:270:VAL:HG23	1:D:271:ASP:N	2.28	0.48
1:A:209:LEU:HD11	1:A:314:LEU:HD11	1.94	0.48
1:A:381:VAL:HG21	1:A:391:VAL:CG2	2.44	0.48
1:A:300:LYS:HE2	2:E:2:DG:O6	2.14	0.48
2:G:2:DG:H2"	2:G:3:DG:OP1	2.13	0.48
1:B:225:ARG:CG	1:C:282:TYR:CZ	2.97	0.47
1:D:248:LYS:HB3	1:D:270:VAL:HG21	1.97	0.47
1:A:422:ILE:HG23	1:A:426:LEU:HD23	1.97	0.47
1:B:423:PHE:CD1	1:B:457:LYS:HE3	2.49	0.47
1:A:205:LEU:HD23	1:A:206:SER:N	2.30	0.47
1:A:423:PHE:HB3	1:A:424:PRO:CD	2.41	0.47
1:B:401:SER:OG	1:B:402:ASP:N	2.47	0.47
1:B:423:PHE:HB3	1:B:424:PRO:CD	2.45	0.47
1:C:205:LEU:HD23	1:C:206:SER:H	1.80	0.47
1:A:392:LEU:HD13	1:A:441:TYR:CD1	2.50	0.47
1:C:233:ARG:HD3	1:C:282:TYR:OH	2.14	0.47
1:C:384:THR:CG2	1:C:386:LYS:HG2	2.44	0.47
1:D:451:ASN:O	1:D:452:ASP:HB2	2.15	0.47
1:D:293:PHE:HB2	1:D:298:GLU:HB3	1.97	0.47
1:D:392:LEU:HD13	1:D:441:TYR:CE1	2.49	0.47
1:C:449:ILE:HG13	1:C:454:VAL:HG23	1.97	0.46
1:B:403:LYS:O	1:B:424:PRO:HD3	2.15	0.46
1:C:378:ILE:HD13	1:C:379:SER:N	2.28	0.46
1:A:269:LYS:HE2	1:A:304:GLY:HA2	1.98	0.46
1:A:443:ASN:HB2	1:A:463:SER:OG	2.16	0.46
1:D:209:LEU:HD11	1:D:314:LEU:HD11	1.97	0.46
1:B:399:GLN:HB2	1:B:406:ARG:HB2	1.98	0.46
1:B:312:GLU:O	1:B:313:THR:HB	2.15	0.46
1:B:300:LYS:HE2	2:F:2:DG:O6	2.15	0.46
1:A:278:LYS:HE3	1:A:460:ILE:O	2.16	0.46
1:B:422:ILE:HG23	1:B:426:LEU:HD23	1.98	0.46
1:C:231:GLU:OE1	1:C:311:LYS:NZ	2.48	0.46
1:B:207:ASP:O	1:B:210:SER:HB2	2.16	0.45
1:C:419:LYS:HE3	1:C:419:LYS:HB2	1.77	0.45
1:D:424:PRO:O	1:D:425:ASP:CB	2.63	0.45
1:D:264:LEU:HG	1:D:299:LEU:HB2	1.99	0.45
1:A:423:PHE:CB	1:A:424:PRO:HD2	2.44	0.45
1:B:216:ILE:HD13	1:B:262:ILE:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:PHE:CB	1:A:424:PRO:CD	2.95	0.45
1:B:225:ARG:H	1:B:225:ARG:HG3	1.64	0.45
1:B:390:SER:N	3:B:2:HOH:O	2.49	0.45
1:C:449:ILE:HG13	1:C:454:VAL:CG2	2.47	0.45
1:D:458:ASN:HB2	1:D:462:SER:HB2	1.99	0.45
1:D:379:SER:O	1:D:382:LEU:CD2	2.65	0.45
1:A:214:LEU:HD11	1:A:223:LYS:HE2	1.98	0.45
1:D:311:LYS:O	1:D:314:LEU:HB2	2.17	0.44
1:C:447:ARG:CZ	1:C:454:VAL:HG21	2.48	0.44
1:A:311:LYS:O	1:A:314:LEU:HB2	2.17	0.44
1:A:450:TYR:CD2	1:A:450:TYR:C	2.91	0.44
1:B:399:GLN:HG3	3:B:14:HOH:O	2.17	0.44
1:B:448:GLN:NE2	1:B:457:LYS:HD3	2.32	0.43
1:C:214:LEU:HD11	1:C:223:LYS:HE3	1.99	0.43
1:A:294:LYS:C	1:A:296:ASN:H	2.21	0.43
1:A:390:SER:N	3:A:3:HOH:O	2.51	0.43
1:A:429:GLU:HG3	1:A:430:TRP:CD1	2.52	0.43
1:D:378:ILE:CG1	1:D:411:ASP:HB2	2.48	0.43
1:D:394:GLN:HA	1:D:438:LYS:O	2.18	0.43
1:A:297:LEU:HA	1:A:297:LEU:HD12	1.79	0.43
1:D:378:ILE:HG12	1:D:411:ASP:CG	2.38	0.43
1:A:424:PRO:O	1:A:425:ASP:HB2	2.18	0.43
1:B:411:ASP:OD1	1:B:411:ASP:N	2.51	0.43
1:B:423:PHE:CE1	1:B:457:LYS:HE3	2.53	0.43
1:D:375:THR:HA	1:D:390:SER:O	2.18	0.43
1:B:265:SER:HB2	1:B:267:TRP:NE1	2.33	0.43
1:B:311:LYS:C	1:B:312:GLU:O	2.57	0.43
1:C:233:ARG:HB2	1:C:282:TYR:CE1	2.53	0.43
1:C:392:LEU:HD11	1:C:467:SER:HB2	2.01	0.43
1:B:294:LYS:C	1:B:296:ASN:H	2.20	0.43
1:C:300:LYS:HE2	1:C:302:LYS:HD2	2.00	0.43
1:C:403:LYS:HG2	1:C:428:TYR:OH	2.19	0.43
1:B:390:SER:HA	1:B:442:PHE:O	2.19	0.42
1:D:379:SER:HB3	1:D:414:CYS:SG	2.59	0.42
1:D:435:SER:HB2	1:D:438:LYS:HD2	2.00	0.42
1:A:248:LYS:O	1:A:270:VAL:HG21	2.19	0.42
1:B:294:LYS:O	1:B:295:ASN:HB2	2.18	0.42
1:A:244:GLN:HG3	1:B:387:GLN:OE1	2.19	0.42
1:C:451:ASN:C	1:C:453:GLU:H	2.23	0.42
1:D:396:VAL:HG11	1:D:410:CYS:SG	2.60	0.42
1:C:378:ILE:N	1:C:378:ILE:HD13	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:293:PHE:HB2	1:D:298:GLU:HG2	2.02	0.42
1:D:423:PHE:HB3	1:D:424:PRO:CD	2.45	0.42
1:A:435:SER:HB2	1:A:438:LYS:HD2	2.02	0.42
1:A:378:ILE:HD13	1:A:411:ASP:HB3	2.01	0.41
1:B:438:LYS:HB3	1:B:440:TYR:CE1	2.54	0.41
1:B:390:SER:HB3	1:B:443:ASN:HA	2.01	0.41
1:D:289:GLU:OE2	1:D:303:PHE:CE1	2.73	0.41
1:C:378:ILE:HD11	1:C:414:CYS:H	1.85	0.41
1:D:375:THR:C	1:D:376:LEU:HD12	2.41	0.41
1:A:377:LEU:O	1:A:381:VAL:HG23	2.19	0.41
1:A:389:LEU:HD21	1:A:447:ARG:HG3	2.02	0.41
1:A:234:VAL:HG22	1:A:252:ILE:HD12	2.03	0.41
1:D:297:LEU:HA	1:D:297:LEU:HD12	1.68	0.41
1:C:379:SER:O	1:C:382:LEU:HB3	2.21	0.41
1:C:378:ILE:CG1	1:C:414:CYS:HB3	2.50	0.41
1:B:250:PHE:HD2	1:B:270:VAL:HG13	1.85	0.41
1:A:375:THR:O	1:A:375:THR:HG23	2.20	0.40
1:B:225:ARG:HD2	1:C:282:TYR:CE1	2.56	0.40
1:A:451:ASN:C	1:A:453:GLU:N	2.73	0.40
1:C:424:PRO:O	1:C:425:ASP:CB	2.66	0.40
1:D:380:GLU:O	1:D:383:LYS:N	2.52	0.40
1:D:411:ASP:HB3	1:D:413:SER:H	1.85	0.40
1:D:455:GLN:O	1:D:455:GLN:HG3	2.20	0.40
1:B:406:ARG:NH1	1:B:419:LYS:HZ3	2.07	0.40
1:C:408:LYS:HA	1:C:418:LEU:O	2.22	0.40
1:A:376:LEU:HD12	1:A:390:SER:O	2.22	0.40
1:C:443:ASN:HB2	1:C:463:SER:OG	2.21	0.40
1:B:378:ILE:HB	1:B:411:ASP:CG	2.42	0.40
1:B:406:ARG:NH1	1:B:419:LYS:HZ1	2.16	0.40
1:C:384:THR:HG22	1:C:386:LYS:HG2	2.03	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	211/213 (99%)	193 (92%)	15 (7%)	3 (1%)	11	21
1	B	208/213 (98%)	189 (91%)	18 (9%)	1 (0%)	29	50
1	C	209/213 (98%)	195 (93%)	13 (6%)	1 (0%)	29	50
1	D	208/213 (98%)	193 (93%)	12 (6%)	3 (1%)	11	21
All	All	836/852 (98%)	770 (92%)	58 (7%)	8 (1%)	15	30

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	425	ASP
1	B	425	ASP
1	C	425	ASP
1	D	425	ASP
1	A	295	ASN
1	D	294	LYS
1	A	294	LYS
1	D	424	PRO

### 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	199/199 (100%)	176 (88%)	23 (12%)	5	9
1	B	197/199 (99%)	172 (87%)	25 (13%)	4	7
1	C	197/199 (99%)	179 (91%)	18 (9%)	9	17
1	D	197/199 (99%)	185 (94%)	12 (6%)	18	36
All	All	790/796 (99%)	712 (90%)	78 (10%)	8	14

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	205	LEU
1	A	210	SER
1	A	221	VAL
1	A	233	ARG
1	A	239	SER
1	A	268	GLN
1	A	297	LEU
1	A	302	LYS
1	A	320	SER
1	A	376	LEU
1	A	382	LEU
1	A	389	LEU
1	A	402	ASP
1	A	414	CYS
1	A	417	GLU
1	A	418	LEU
1	A	429	GLU
1	A	431	ARG
1	A	432	ASP
1	A	447	ARG
1	A	449	ILE
1	A	451	ASN
1	A	466	GLU
1	B	205	LEU
1	B	207	ASP
1	B	225	ARG
1	B	244	GLN
1	B	252	ILE
1	B	264	LEU
1	B	293	PHE
1	B	376	LEU
1	B	379	SER
1	B	380	GLU
1	B	382	LEU
1	B	386	LYS
1	B	389	LEU
1	B	402	ASP
1	B	408	LYS
1	B	410	CYS
1	B	414	CYS
1	B	417	GLU
1	B	418	LEU
1	B	438	LYS

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Mol	Chain	Res	Type
1	B	449	ILE
1	B	450	TYR
1	B	453	GLU
1	B	454	VAL
1	B	466	GLU
1	C	221	VAL
1	C	223	LYS
1	C	225	ARG
1	C	233	ARG
1	C	276	THR
1	C	281	GLN
1	C	313	THR
1	C	314	LEU
1	C	377	LEU
1	C	378	ILE
1	C	384	THR
1	C	401	SER
1	C	412	ASN
1	C	417	GLU
1	C	418	LEU
1	C	429	GLU
1	C	432	ASP
1	C	449	ILE
1	D	233	ARG
1	D	248	LYS
1	D	268	GLN
1	D	275	GLN
1	D	281	GLN
1	D	297	LEU
1	D	311	LYS
1	D	320	SER
1	D	403	LYS
1	D	432	ASP
1	D	449	ILE
1	D	450	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	296	ASN
1	B	448	GLN
1	C	461	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	213/213 (100%)	0.69	22 (10%) <b>6</b> <b>4</b>	32, 54, 106, 124	0
1	B	210/213 (98%)	0.60	15 (7%) <b>16</b> <b>11</b>	32, 54, 103, 125	0
1	C	211/213 (99%)	0.63	20 (9%) <b>8</b> <b>5</b>	33, 54, 105, 122	0
1	D	210/213 (98%)	0.61	17 (8%) <b>12</b> <b>8</b>	32, 54, 103, 129	0
2	E	4/4 (100%)	0.16	0 <b>100</b> <b>100</b>	48, 51, 58, 70	0
2	F	4/4 (100%)	0.30	0 <b>100</b> <b>100</b>	49, 51, 60, 75	0
2	G	4/4 (100%)	0.50	0 <b>100</b> <b>100</b>	48, 50, 59, 70	0
2	H	4/4 (100%)	0.13	0 <b>100</b> <b>100</b>	49, 49, 60, 74	0
All	All	860/868 (99%)	0.63	74 (8%) <b>10</b> <b>7</b>	32, 54, 104, 129	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	411	ASP	13.1
1	D	415	ASN	9.0
1	A	415	ASN	6.8
1	B	454	VAL	6.6
1	A	392	LEU	6.2
1	B	452	ASP	5.9
1	C	411	ASP	5.6
1	D	378	ILE	5.5
1	A	412	ASN	5.3
1	C	452	ASP	5.0
1	D	294	LYS	4.5
1	A	452	ASP	4.4
1	D	413	SER	4.4
1	B	451	ASN	4.4
1	C	415	ASN	4.3
1	B	247	SER	4.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	450	TYR	4.2
1	C	376	LEU	4.1
1	D	246	ASP	4.1
1	D	432	ASP	4.1
1	D	377	LEU	4.1
1	A	414	CYS	4.0
1	B	382	LEU	4.0
1	D	379	SER	3.9
1	D	454	VAL	3.9
1	A	413	SER	3.8
1	C	412	ASN	3.8
1	D	414	CYS	3.8
1	A	375	THR	3.7
1	C	454	VAL	3.6
1	B	431	ARG	3.5
1	B	450	TYR	3.5
1	C	319	GLY	3.5
1	A	384	THR	3.3
1	D	247	SER	3.3
1	C	318	GLY	3.2
1	A	439	TRP	3.1
1	D	411	ASP	3.1
1	A	402	ASP	3.1
1	B	313	THR	3.1
1	A	293	PHE	3.0
1	B	246	ASP	2.9
1	D	403	LYS	2.9
1	A	377	LEU	2.8
1	C	389	LEU	2.8
1	C	383	LYS	2.8
1	C	380	GLU	2.6
1	D	293	PHE	2.6
1	C	375	THR	2.6
1	C	384	THR	2.6
1	C	385	SER	2.6
1	A	245	GLN	2.5
1	A	453	GLU	2.5
1	A	223	LYS	2.5
1	A	246	ASP	2.5
1	A	383	LYS	2.4
1	C	403	LYS	2.4
1	A	432	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	381	VAL	2.4
1	C	447	ARG	2.4
1	D	417	GLU	2.3
1	D	425	ASP	2.3
1	C	245	GLN	2.3
1	C	382	LEU	2.3
1	B	380	GLU	2.2
1	A	389	LEU	2.2
1	A	437	ASN	2.2
1	B	403	LYS	2.1
1	B	402	ASP	2.1
1	B	447	ARG	2.1
1	B	293	PHE	2.1
1	C	314	LEU	2.1
1	B	377	LEU	2.0
1	D	389	LEU	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.