



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:29 PM GMT

PDB ID : 1XBT  
Title : Crystal Structure of Human Thymidine Kinase 1  
Authors : Welin, M.; Kosinska, U.; Mikkelsen, N.E.; Carnrot, C.; Zhu, C.; Wang, L.; Eriksson, S.; Munch-Petersen, B.; Eklund, H.  
Deposited on : 2004-08-31  
Resolution : 2.40 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  
<http://wwpdb.org/validation/2016/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  
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

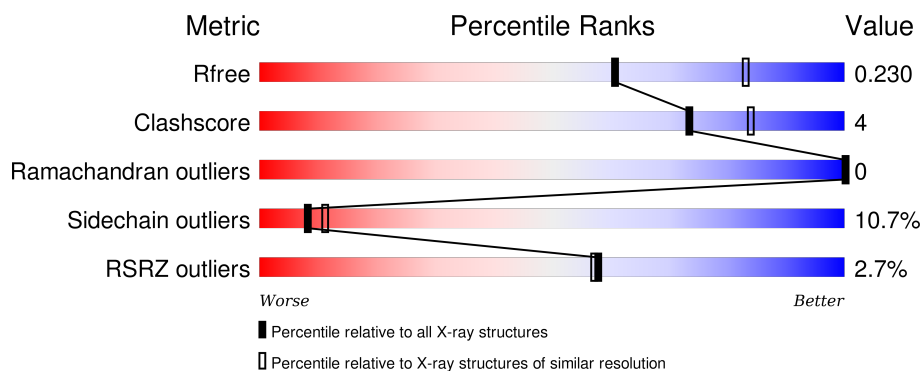
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	193	
1	B	193	
1	C	193	
1	D	193	
1	E	193	

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Mol	Chain	Length	Quality of chain
1	F	193	<div><div><div></div><div></div><div></div><div></div></div><div><div>4%</div><div>69%</div><div>12%</div><div>•</div><div>17%</div></div></div>
1	G	193	<div><div><div></div><div></div><div></div><div></div></div><div><div>2%</div><div>72%</div><div>12%</div><div>•</div><div>15%</div></div></div>
1	H	193	<div><div><div></div><div></div><div></div><div></div></div><div><div>2%</div><div>67%</div><div>14%</div><div>•</div><div>17%</div></div></div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10451 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 Thymidine kinase, cytosolic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	160	Total	C	N	O	S	0	0	0
			1239	794	215	219	11			
1	B	160	Total	C	N	O	S	0	0	0
			1239	794	215	219	11			
1	C	160	Total	C	N	O	S	0	0	0
			1239	794	215	219	11			
1	D	162	Total	C	N	O	S	0	0	0
			1257	806	217	223	11			
1	E	160	Total	C	N	O	S	0	0	0
			1239	794	215	219	11			
1	F	161	Total	C	N	O	S	0	0	0
			1248	800	217	220	11			
1	G	165	Total	C	N	O	S	0	0	0
			1278	817	220	230	11			
1	H	161	Total	C	N	O	S	0	0	0
			1251	803	216	221	11			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	H	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

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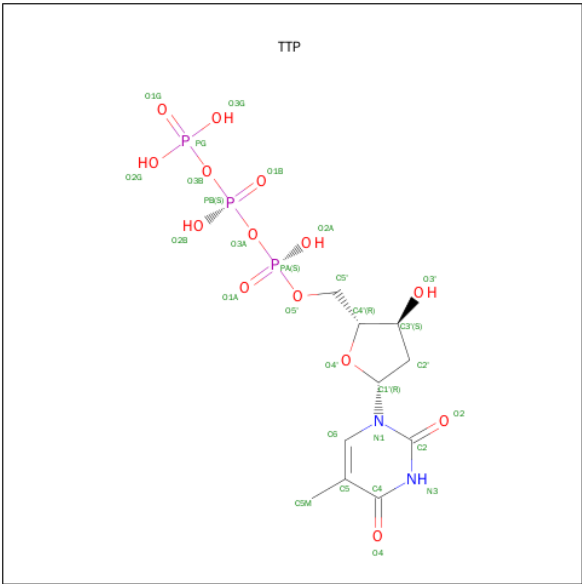
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	H	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is THYMIDINE-5'-TRIPHOSPHATE (three-letter code: TTP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>2</sub>O<sub>14</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	B	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	C	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	D	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	E	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	F	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	G	1	Total	C	N	O	P	0	0
			29	10	2	14	3		
4	H	1	Total	C	N	O	P	0	0
			29	10	2	14	3		

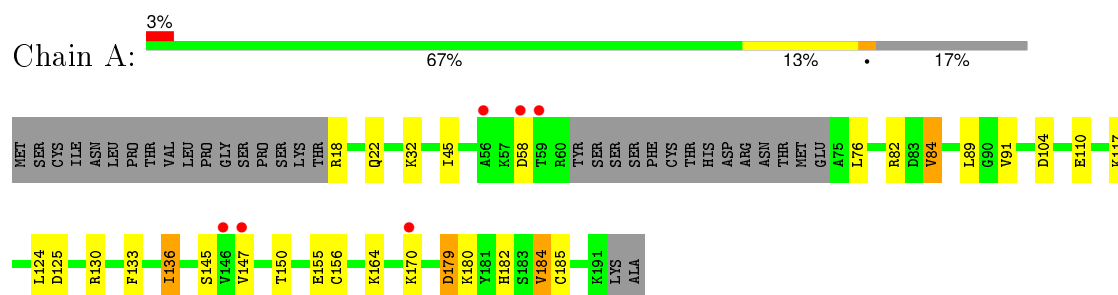
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	23	Total	O	0	0
			23	23		
5	B	31	Total	O	0	0
			31	31		
5	C	20	Total	O	0	0
			20	20		
5	D	35	Total	O	0	0
			35	35		
5	E	24	Total	O	0	0
			24	24		
5	F	25	Total	O	0	0
			25	25		
5	G	33	Total	O	0	0
			33	33		
5	H	22	Total	O	0	0
			22	22		

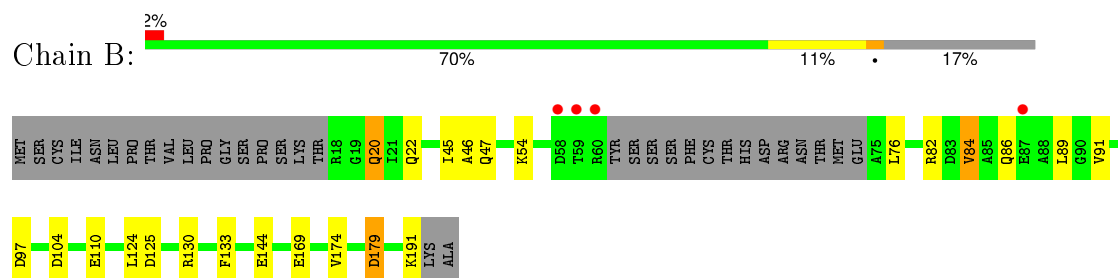
### 3 Residue-property plots [i](#)

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.

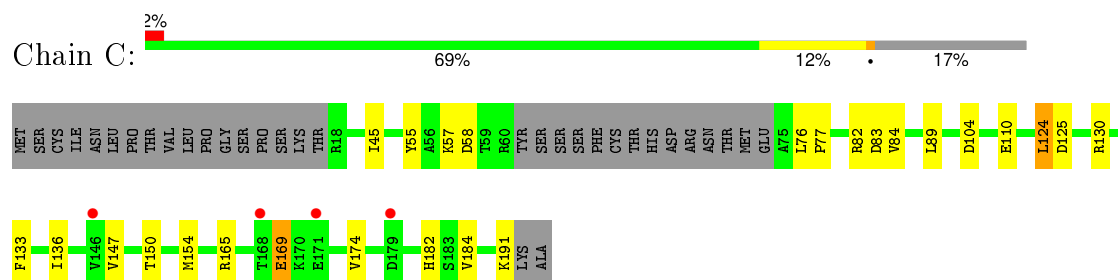
- Molecule 1: Thymidine kinase, cytosolic



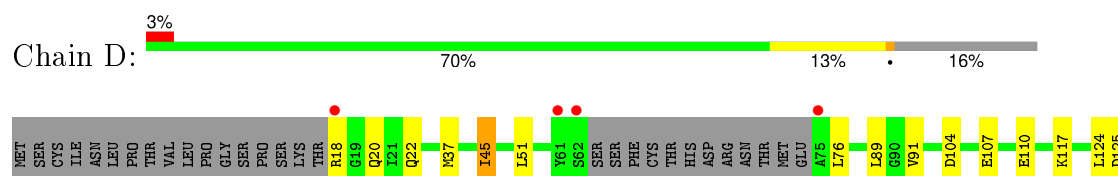
- Molecule 1: Thymidine kinase, cytosolic

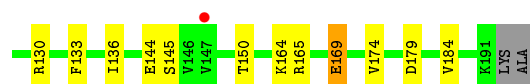


- Molecule 1: Thymidine kinase, cytosolic



- Molecule 1: Thymidine kinase, cytosolic





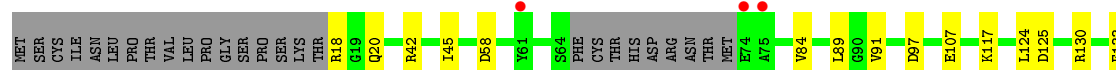
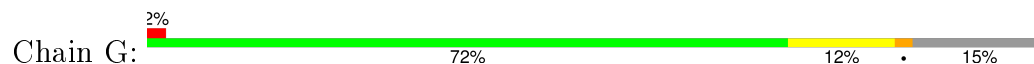
- Molecule 1: Thymidine kinase, cytosolic



- Molecule 1: Thymidine kinase, cytosolic



- Molecule 1: Thymidine kinase, cytosolic



- Molecule 1: Thymidine kinase, cytosolic





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.67Å 123.18Å 116.00Å 90.00° 130.20° 90.00°	Depositor
Resolution (Å)	42.07 – 2.40 42.07 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.9 (42.07-2.40) 98.9 (42.07-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.17 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.201 , 0.232 0.201 , 0.230	Depositor DCC
$R_{free}$ test set	3317 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.8	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 42.9	EDS
Estimated twinning fraction	0.010 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 65407 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10451	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, TTP, 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	0.45	0/1257	0.69	4/1690 (0.2%)
1	B	0.46	0/1257	0.70	4/1690 (0.2%)
1	C	0.46	0/1257	0.70	4/1690 (0.2%)
1	D	0.46	0/1276	0.71	3/1716 (0.2%)
1	E	0.45	0/1257	0.71	4/1690 (0.2%)
1	F	0.46	0/1266	0.69	2/1701 (0.1%)
1	G	0.46	0/1297	0.72	4/1744 (0.2%)
1	H	0.43	0/1270	0.70	3/1708 (0.2%)
All	All	0.45	0/10137	0.70	28/13629 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	104	ASP	CB-CG-OD2	6.97	124.57	118.30
1	B	104	ASP	CB-CG-OD2	6.70	124.33	118.30
1	A	104	ASP	CB-CG-OD2	6.51	124.16	118.30
1	D	104	ASP	CB-CG-OD2	6.49	124.14	118.30
1	C	104	ASP	CB-CG-OD2	6.47	124.12	118.30
1	G	97	ASP	CB-CG-OD2	6.38	124.04	118.30
1	E	179	ASP	CB-CG-OD2	6.18	123.86	118.30
1	C	125	ASP	CB-CG-OD2	6.13	123.82	118.30
1	B	125	ASP	CB-CG-OD2	6.00	123.70	118.30
1	H	58	ASP	CB-CG-OD2	5.90	123.61	118.30
1	G	125	ASP	CB-CG-OD2	5.87	123.58	118.30
1	A	179	ASP	CB-CG-OD2	5.73	123.46	118.30
1	B	179	ASP	CB-CG-OD2	5.71	123.44	118.30
1	G	58	ASP	CB-CG-OD2	5.68	123.41	118.30
1	E	97	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	58	ASP	CB-CG-OD2	5.36	123.12	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	58	ASP	CB-CG-OD2	5.28	123.05	118.30
1	D	125	ASP	CB-CG-OD2	5.27	123.04	118.30
1	G	179	ASP	CB-CG-OD2	5.26	123.04	118.30
1	A	125	ASP	CB-CG-OD2	5.25	123.03	118.30
1	C	83	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	97	ASP	CB-CG-OD2	5.22	123.00	118.30
1	H	97	ASP	CB-CG-OD2	5.22	123.00	118.30
1	D	179	ASP	CB-CG-OD2	5.20	122.98	118.30
1	E	125	ASP	CB-CG-OD2	5.11	122.90	118.30
1	F	58	ASP	CB-CG-OD2	5.07	122.86	118.30
1	F	125	ASP	CB-CG-OD2	5.02	122.81	118.30
1	H	179	ASP	CB-CG-OD2	5.00	122.80	118.30

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	1239	0	1280	9	0
1	B	1239	0	1280	12	0
1	C	1239	0	1280	11	0
1	D	1257	0	1293	11	0
1	E	1239	0	1279	13	0
1	F	1248	0	1293	13	0
1	G	1278	0	1309	14	0
1	H	1251	0	1289	14	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	29	0	13	0	0
4	B	29	0	13	1	0
4	C	29	0	13	2	0
4	D	29	0	13	0	0
4	E	29	0	13	0	0
4	F	29	0	13	0	0
4	G	29	0	13	0	0
4	H	29	0	13	0	0
5	A	23	0	0	0	0
5	B	31	0	0	2	0
5	C	20	0	0	0	0
5	D	35	0	0	0	0
5	E	24	0	0	0	0
5	F	25	0	0	0	0
5	G	33	0	0	0	0
5	H	22	0	0	1	0
All	All	10451	0	10407	84	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 4.

All (84) 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:D:107:GLU:HG3	1:G:107:GLU:HG3	1.40	1.03
1:B:20:GLN:HG2	1:B:144:GLU:HG3	1.53	0.90
1:F:158:ARG:HH21	1:G:158:ARG:HH21	1.28	0.81
1:F:20:GLN:HG2	1:F:144:GLU:HG3	1.64	0.79
1:A:76:LEU:HD13	1:A:84:VAL:HG22	1.73	0.71
1:D:20:GLN:HG2	1:D:144:GLU:HG3	1.72	0.69
1:H:82:ARG:HD3	5:H:8199:HOH:O	1.91	0.69
1:F:20:GLN:HG2	1:F:144:GLU:CG	2.23	0.68
1:E:110:GLU:OE1	1:G:130:ARG:NH2	2.26	0.68
1:A:110:GLU:OE1	1:C:130:ARG:NH2	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:158:ARG:NH2	1:G:158:ARG:HH21	1.92	0.66
1:B:130:ARG:NH2	1:D:110:GLU:OE1	2.25	0.64
1:B:110:GLU:OE1	1:D:130:ARG:NH2	2.31	0.63
1:E:20:GLN:HG3	1:E:144:GLU:HG3	1.82	0.61
1:E:20:GLN:CG	1:E:144:GLU:HG3	2.32	0.60
1:B:76:LEU:HD13	1:B:84:VAL:HG22	1.84	0.59
1:D:20:GLN:HG2	1:D:144:GLU:CG	2.32	0.59
1:F:165:ARG:HD3	1:F:169:GLU:HG2	1.85	0.58
1:B:20:GLN:HG2	1:B:144:GLU:CG	2.30	0.58
1:E:37:MET:CE	1:E:52:VAL:HG11	2.34	0.57
4:B:2195:TTP:PA	5:B:2226:HOH:O	2.61	0.57
1:C:84:VAL:CG1	1:C:84:VAL:O	2.52	0.57
1:H:169:GLU:HG3	1:H:174:VAL:HG21	1.88	0.56
1:H:20:GLN:HG2	1:H:144:GLU:HG3	1.86	0.56
1:E:20:GLN:NE2	1:E:22:GLN:HE21	2.05	0.55
1:C:169:GLU:HG3	1:C:174:VAL:HG21	1.89	0.54
1:G:169:GLU:HG3	1:G:174:VAL:HG21	1.89	0.54
1:C:84:VAL:HG12	1:C:84:VAL:O	2.07	0.53
1:E:110:GLU:CD	1:G:130:ARG:HH22	2.12	0.52
1:B:169:GLU:HG3	1:B:174:VAL:HG21	1.92	0.52
1:B:110:GLU:CD	1:D:130:ARG:HH22	2.13	0.52
1:E:37:MET:HE2	1:E:52:VAL:HG11	1.92	0.51
1:A:22:GLN:HG2	1:A:145:SER:HB3	1.93	0.51
1:G:165:ARG:HD3	1:G:169:GLU:HG2	1.93	0.51
1:C:124:LEU:HD21	4:C:3195:TTP:H5'2	1.93	0.50
1:F:191:LYS:HE3	1:F:192:LYS:HZ2	1.76	0.50
1:G:184:VAL:HG13	1:G:185:CYS:O	2.12	0.50
1:D:169:GLU:HG3	1:D:174:VAL:HG21	1.93	0.50
1:A:76:LEU:HD13	1:A:84:VAL:CG2	2.40	0.49
1:F:185:CYS:SG	1:H:18:ARG:NH2	2.85	0.49
1:G:20:GLN:HB2	1:G:144:GLU:CG	2.42	0.49
1:G:155:GLU:OE1	1:G:182:HIS:HD2	1.93	0.49
1:D:165:ARG:HD3	1:D:169:GLU:HG2	1.95	0.48
1:G:84:VAL:CG1	1:G:84:VAL:O	2.61	0.48
1:G:42:ARG:HH21	1:H:150:THR:HG22	1.78	0.48
1:F:169:GLU:HG3	1:F:174:VAL:HG21	1.96	0.47
1:H:55:TYR:CE2	1:H:57:LYS:HB2	2.49	0.47
1:D:22:GLN:HG2	1:D:145:SER:HB3	1.97	0.47
1:B:20:GLN:NE2	1:B:22:GLN:HE21	2.13	0.47
1:F:37:MET:CE	1:F:52:VAL:HG11	2.46	0.46
1:B:76:LEU:HD13	1:B:84:VAL:CG2	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:20:GLN:HG2	1:H:144:GLU:CG	2.45	0.46
1:C:55:TYR:CE2	1:C:57:LYS:HB2	2.50	0.46
1:F:130:ARG:NH2	1:H:110:GLU:OE1	2.46	0.46
1:F:76:LEU:HD13	1:F:84:VAL:CG2	2.46	0.45
1:E:37:MET:HE1	1:E:52:VAL:HG11	1.98	0.45
1:H:184:VAL:HG13	1:H:185:CYS:O	2.17	0.45
1:H:136:ILE:HD13	1:H:139:LEU:HD12	1.99	0.45
1:A:32:LYS:HE2	1:A:32:LYS:HB2	1.83	0.44
1:B:20:GLN:HE22	1:B:22:GLN:HE21	1.66	0.44
1:E:20:GLN:HG2	1:E:144:GLU:HG3	2.00	0.44
1:E:136:ILE:HD13	1:E:136:ILE:O	2.17	0.43
1:C:76:LEU:HA	1:C:77:PRO:HD3	1.92	0.43
1:G:20:GLN:HB2	1:G:144:GLU:HG3	2.01	0.43
1:E:55:TYR:CE2	1:E:57:LYS:HB2	2.54	0.43
1:C:165:ARG:HD3	1:C:169:GLU:HG2	2.00	0.43
1:A:184:VAL:HG13	1:A:185:CYS:O	2.19	0.42
1:F:191:LYS:C	1:F:192:LYS:HG2	2.40	0.42
1:E:165:ARG:HD3	1:E:169:GLU:HG2	2.01	0.42
1:H:20:GLN:NE2	1:H:22:GLN:HE21	2.17	0.42
1:H:165:ARG:HD3	1:H:169:GLU:HG2	2.02	0.41
1:E:20:GLN:HG2	1:E:144:GLU:CG	2.49	0.41
1:A:155:GLU:OE1	1:A:182:HIS:HD2	2.03	0.41
1:A:136:ILE:HD13	1:A:136:ILE:O	2.20	0.41
1:G:107:GLU:CD	1:G:107:GLU:H	2.22	0.41
1:B:46:ALA:O	1:B:47:GLN:HB2	2.20	0.41
1:A:130:ARG:NH2	1:C:110:GLU:OE1	2.46	0.41
1:D:51:LEU:HD11	1:D:76:LEU:HD12	2.03	0.41
1:H:20:GLN:HE22	1:H:22:GLN:HE21	1.69	0.41
1:C:124:LEU:CD2	4:C:3195:TTP:H5'2	2.51	0.40
1:C:154:MET:CE	1:D:45:ILE:HG23	2.51	0.40
1:H:84:VAL:O	1:H:84:VAL:HG13	2.21	0.40
1:B:130:ARG:NH2	5:B:2212:HOH:O	2.41	0.40
1:F:37:MET:HE1	1:F:52:VAL:HG11	2.03	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	156/193 (81%)	156 (100%)	0	0	100	100
1	B	156/193 (81%)	154 (99%)	2 (1%)	0	100	100
1	C	156/193 (81%)	155 (99%)	1 (1%)	0	100	100
1	D	158/193 (82%)	158 (100%)	0	0	100	100
1	E	156/193 (81%)	155 (99%)	1 (1%)	0	100	100
1	F	157/193 (81%)	156 (99%)	1 (1%)	0	100	100
1	G	161/193 (83%)	160 (99%)	1 (1%)	0	100	100
1	H	157/193 (81%)	156 (99%)	1 (1%)	0	100	100
All	All	1257/1544 (81%)	1250 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	130/161 (81%)	112 (86%)	18 (14%)	4	5
1	B	130/161 (81%)	118 (91%)	12 (9%)	11	16
1	C	130/161 (81%)	118 (91%)	12 (9%)	11	16
1	D	132/161 (82%)	119 (90%)	13 (10%)	10	14
1	E	130/161 (81%)	114 (88%)	16 (12%)	6	7
1	F	131/161 (81%)	117 (89%)	14 (11%)	8	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	135/161 (84%)	123 (91%)	12 (9%)	12	18
1	H	131/161 (81%)	116 (88%)	15 (12%)	7	9
All	All	1049/1288 (81%)	937 (89%)	112 (11%)	8	11

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	45	ILE
1	A	82	ARG
1	A	84	VAL
1	A	89	LEU
1	A	91	VAL
1	A	117	LYS
1	A	124	LEU
1	A	133	PHE
1	A	136	ILE
1	A	147	VAL
1	A	150	THR
1	A	156	CYS
1	A	164	LYS
1	A	170	LYS
1	A	179	ASP
1	A	180	LYS
1	A	184	VAL
1	B	20	GLN
1	B	45	ILE
1	B	54	LYS
1	B	82	ARG
1	B	84	VAL
1	B	86	GLN
1	B	89	LEU
1	B	91	VAL
1	B	124	LEU
1	B	133	PHE
1	B	179	ASP
1	B	191	LYS
1	C	45	ILE
1	C	82	ARG
1	C	89	LEU
1	C	124	LEU

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Mol	Chain	Res	Type
1	C	133	PHE
1	C	136	ILE
1	C	147	VAL
1	C	150	THR
1	C	169	GLU
1	C	182	HIS
1	C	184	VAL
1	C	191	LYS
1	D	18	ARG
1	D	37	MET
1	D	45	ILE
1	D	89	LEU
1	D	91	VAL
1	D	117	LYS
1	D	124	LEU
1	D	133	PHE
1	D	136	ILE
1	D	150	THR
1	D	164	LYS
1	D	169	GLU
1	D	184	VAL
1	E	18	ARG
1	E	45	ILE
1	E	54	LYS
1	E	82	ARG
1	E	84	VAL
1	E	89	LEU
1	E	91	VAL
1	E	124	LEU
1	E	133	PHE
1	E	136	ILE
1	E	150	THR
1	E	170	LYS
1	E	171	GLU
1	E	179	ASP
1	E	184	VAL
1	E	191	LYS
1	F	18	ARG
1	F	20	GLN
1	F	45	ILE
1	F	49	LYS
1	F	76	LEU

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Mol	Chain	Res	Type
1	F	89	LEU
1	F	117	LYS
1	F	124	LEU
1	F	133	PHE
1	F	136	ILE
1	F	147	VAL
1	F	150	THR
1	F	180	LYS
1	F	192	LYS
1	G	18	ARG
1	G	45	ILE
1	G	89	LEU
1	G	91	VAL
1	G	117	LYS
1	G	124	LEU
1	G	133	PHE
1	G	136	ILE
1	G	144	GLU
1	G	147	VAL
1	G	169	GLU
1	G	184	VAL
1	H	49	LYS
1	H	59	THR
1	H	79	CYS
1	H	84	VAL
1	H	89	LEU
1	H	117	LYS
1	H	124	LEU
1	H	133	PHE
1	H	136	ILE
1	H	147	VAL
1	H	150	THR
1	H	175	ILE
1	H	180	LYS
1	H	184	VAL
1	H	191	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	182	HIS

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Mol	Chain	Res	Type
1	B	20	GLN
1	B	86	GLN
1	C	20	GLN
1	C	182	HIS
1	E	20	GLN
1	E	129	GLN
1	E	182	HIS
1	G	20	GLN
1	G	86	GLN
1	G	182	HIS
1	H	20	GLN

### 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](#)

Of 24 ligands modelled in this entry, 16 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	TTP	A	1195	3	21,30,30	0.55	0	31,47,47	2.09	3 (9%)
4	TTP	B	2195	3	21,30,30	0.57	0	31,47,47	1.95	3 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	TTP	C	3195	3	21,30,30	0.55	0	31,47,47	1.94	4 (12%)
4	TTP	D	4195	3	21,30,30	0.64	0	31,47,47	2.05	3 (9%)
4	TTP	E	5195	3	21,30,30	0.57	0	31,47,47	2.09	4 (12%)
4	TTP	F	6195	3	21,30,30	0.56	0	31,47,47	1.99	3 (9%)
4	TTP	G	7195	3	21,30,30	0.55	0	31,47,47	1.98	4 (12%)
4	TTP	H	8195	3	21,30,30	0.59	0	31,47,47	1.89	4 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TTP	A	1195	3	-	0/18/34/34	0/2/2/2
4	TTP	B	2195	3	-	0/18/34/34	0/2/2/2
4	TTP	C	3195	3	-	0/18/34/34	0/2/2/2
4	TTP	D	4195	3	-	0/18/34/34	0/2/2/2
4	TTP	E	5195	3	-	0/18/34/34	0/2/2/2
4	TTP	F	6195	3	-	0/18/34/34	0/2/2/2
4	TTP	G	7195	3	-	0/18/34/34	0/2/2/2
4	TTP	H	8195	3	-	0/18/34/34	0/2/2/2

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	4195	TTP	C5-C4-N3	-6.27	118.16	125.14
4	A	1195	TTP	C5-C4-N3	-6.23	118.21	125.14
4	B	2195	TTP	C5-C4-N3	-6.22	118.21	125.14
4	C	3195	TTP	C5-C4-N3	-6.16	118.28	125.14
4	H	8195	TTP	C5-C4-N3	-5.90	118.57	125.14
4	F	6195	TTP	C5-C4-N3	-5.85	118.62	125.14
4	E	5195	TTP	C5-C4-N3	-5.69	118.80	125.14
4	G	7195	TTP	C5-C4-N3	-5.40	119.12	125.14
4	E	5195	TTP	PB-O3A-PA	-4.88	119.03	132.73
4	H	8195	TTP	PB-O3A-PA	-4.00	121.49	132.73
4	D	4195	TTP	PB-O3A-PA	-3.90	121.78	132.73
4	F	6195	TTP	PB-O3A-PA	-3.64	122.49	132.73
4	B	2195	TTP	PB-O3A-PA	-3.59	122.65	132.73
4	C	3195	TTP	PB-O3A-PA	-3.09	124.05	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1195	TTP	PB-O3A-PA	-2.90	124.59	132.73
4	G	7195	TTP	PB-O3A-PA	-2.85	124.72	132.73
4	E	5195	TTP	PB-O3B-PG	-2.60	123.94	132.67
4	G	7195	TTP	O4'-C1'-N1	2.03	111.23	107.72
4	C	3195	TTP	O3G-PG-O1G	2.04	117.15	110.58
4	H	8195	TTP	O3G-PG-O1G	2.07	117.25	110.58
4	H	8195	TTP	C4-N3-C2	6.01	120.44	115.25
4	C	3195	TTP	C4-N3-C2	6.01	120.44	115.25
4	B	2195	TTP	C4-N3-C2	6.08	120.50	115.25
4	E	5195	TTP	C4-N3-C2	6.77	121.10	115.25
4	F	6195	TTP	C4-N3-C2	6.79	121.12	115.25
4	G	7195	TTP	C4-N3-C2	7.22	121.48	115.25
4	D	4195	TTP	C4-N3-C2	7.22	121.48	115.25
4	A	1195	TTP	C4-N3-C2	7.37	121.62	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2195	TTP	1	0
4	C	3195	TTP	2	0

## 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	160/193 (82%)	0.01	6 (3%) 44 45	23, 33, 48, 53	0
1	B	160/193 (82%)	0.10	4 (2%) 61 60	23, 33, 48, 57	0
1	C	160/193 (82%)	0.08	4 (2%) 61 60	23, 33, 48, 54	0
1	D	162/193 (83%)	0.01	5 (3%) 52 52	23, 33, 48, 52	0
1	E	160/193 (82%)	-0.09	1 (0%) 90 90	22, 33, 48, 54	0
1	F	161/193 (83%)	-0.06	8 (4%) 32 33	22, 33, 48, 53	0
1	G	165/193 (85%)	0.05	3 (1%) 71 71	22, 33, 48, 55	0
1	H	161/193 (83%)	-0.00	4 (2%) 61 60	23, 33, 47, 51	0
All	All	1289/1544 (83%)	0.01	35 (2%) 58 57	22, 33, 48, 57	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	59	THR	8.4
1	A	59	THR	4.5
1	G	74	GLU	4.1
1	G	75	ALA	3.9
1	D	75	ALA	3.9
1	E	59	THR	3.5
1	H	75	ALA	3.4
1	F	168	THR	3.0
1	A	170	LYS	2.8
1	B	60	ARG	2.8
1	B	58	ASP	2.6
1	H	86	GLN	2.6
1	A	56	ALA	2.6
1	D	18	ARG	2.6
1	F	18	ARG	2.6
1	D	61	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	171	GLU	2.5
1	F	192	LYS	2.5
1	F	83	ASP	2.4
1	C	168	THR	2.4
1	C	179	ASP	2.2
1	F	191	LYS	2.2
1	F	47	GLN	2.2
1	A	147	VAL	2.2
1	D	147	VAL	2.2
1	A	58	ASP	2.2
1	A	146	VAL	2.1
1	C	146	VAL	2.1
1	D	62	SER	2.1
1	B	87	GLU	2.1
1	H	47	GLN	2.1
1	G	61	TYR	2.0
1	F	75	ALA	2.0
1	H	89	LEU	2.0
1	F	167	GLY	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	TTP	C	3195	29/29	0.93	0.16	0.40	35,37,42,42	0
4	TTP	F	6195	29/29	0.96	0.12	-0.26	26,34,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	TTP	B	2195	29/29	0.95	0.14	-0.29	24,29,39,40	0
4	TTP	A	1195	29/29	0.94	0.13	-0.39	24,31,39,40	0
4	TTP	E	5195	29/29	0.96	0.11	-0.81	26,31,43,44	0
4	TTP	D	4195	29/29	0.98	0.11	-0.95	17,22,31,33	0
2	ZN	G	7193	1/1	0.99	0.08	-1.02	35,35,35,35	0
2	ZN	H	8193	1/1	1.00	0.07	-1.03	35,35,35,35	0
2	ZN	D	4193	1/1	1.00	0.07	-1.10	36,36,36,36	0
4	TTP	H	8195	29/29	0.99	0.10	-1.12	17,22,31,31	0
2	ZN	B	2193	1/1	1.00	0.06	-1.16	30,30,30,30	0
2	ZN	E	5193	1/1	1.00	0.06	-1.32	32,32,32,32	0
2	ZN	F	6193	1/1	0.99	0.05	-1.37	38,38,38,38	0
2	ZN	A	1193	1/1	0.99	0.05	-1.51	31,31,31,31	0
4	TTP	G	7195	29/29	0.98	0.10	-1.76	19,23,31,32	0
2	ZN	C	3193	1/1	0.99	0.06	-1.83	41,41,41,41	0
3	MG	H	8194	1/1	0.89	0.06	-	36,36,36,36	0
3	MG	B	2194	1/1	0.64	0.20	-	41,41,41,41	0
3	MG	C	3194	1/1	0.80	0.10	-	44,44,44,44	0
3	MG	A	1194	1/1	0.51	0.17	-	36,36,36,36	0
3	MG	F	6194	1/1	0.92	0.12	-	48,48,48,48	0
3	MG	G	7194	1/1	0.95	0.06	-	34,34,34,34	0
3	MG	D	4194	1/1	0.90	0.13	-	35,35,35,35	0
3	MG	E	5194	1/1	0.47	0.18	-	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.