



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2023 – 10:28 PM EDT

PDB ID : 6UEQ
Title : Structure of TBP bound to C-C mismatch containing TATA site
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Deposited on : 2019-09-22
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

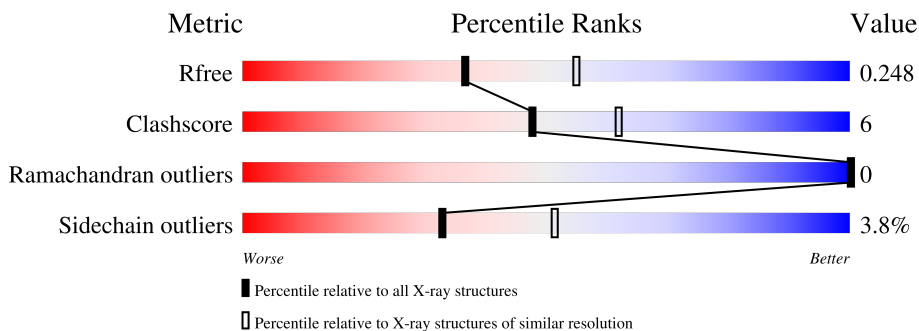
1 Overall quality at a glance

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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	B	219	 73% 11% 15%
2	C	14	 86% 14%
3	D	14	 79% 21%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2120 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 TATA-box-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	187	1470	959	248	255	8	0	0	0

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	MET	-	initiating methionine	UNP P28147
B	-17	SER	-	expression tag	UNP P28147
B	-16	SER	-	expression tag	UNP P28147
B	-15	HIS	-	expression tag	UNP P28147
B	-14	HIS	-	expression tag	UNP P28147
B	-13	HIS	-	expression tag	UNP P28147
B	-12	HIS	-	expression tag	UNP P28147
B	-11	HIS	-	expression tag	UNP P28147
B	-10	HIS	-	expression tag	UNP P28147
B	-9	SER	-	expression tag	UNP P28147
B	-8	SER	-	expression tag	UNP P28147
B	-7	GLY	-	expression tag	UNP P28147
B	-6	LEU	-	expression tag	UNP P28147
B	-5	VAL	-	expression tag	UNP P28147
B	-4	PRO	-	expression tag	UNP P28147
B	-3	ARG	-	expression tag	UNP P28147
B	-2	GLY	-	expression tag	UNP P28147
B	-1	SER	-	expression tag	UNP P28147
B	0	HIS	-	expression tag	UNP P28147

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*CP*TP*AP*TP*AP*AP*AP*AP*CP*GP*GP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	14	286	137	58	78	13	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	D	14	279	136	44	86	13	0	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	B	1	5	4	1	0	0
4	D	1	5	4	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	B	53	53	53	0	0
5	C	9	9	9	0	0
5	D	13	13	13	0	0

4 Data and refinement statistics i

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	45.40Å 45.58Å 155.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.61 – 2.40 77.62 – 2.09	Depositor EDS
% Data completeness (in resolution range)	96.6 (77.61-2.40) 91.0 (77.62-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.08Å)	Xtrriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.212 , 0.249 0.212 , 0.248	Depositor DCC
R_{free} test set	1806 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtrriage
Anisotropy	0.314	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.036 for k,h,-l	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2120	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	B	0.27	0/1500	0.45	0/2020
2	C	0.52	0/322	0.78	0/495
3	D	0.58	0/310	1.06	0/476
All	All	0.38	0/2132	0.64	0/2991

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1470	0	1547	20	0
2	C	286	0	158	1	0
3	D	279	0	162	3	0
4	B	5	0	0	0	0
4	D	5	0	0	1	0
5	B	53	0	0	6	1
5	C	9	0	0	0	0
5	D	13	0	0	1	0
All	All	2120	0	1867	23	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 6.

All (23) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:226:DA:OP2	5:D:401:HOH:O	1.83	0.95
1:B:131:GLU:OE2	5:B:401:HOH:O	1.96	0.83
1:B:64:ILE:O	5:B:402:HOH:O	2.06	0.71
1:B:12:VAL:N	5:B:404:HOH:O	2.29	0.65
1:B:196:ARG:NH2	5:B:407:HOH:O	2.34	0.60
1:B:158:PRO:HD3	1:B:183:LYS:HG2	1.84	0.59
1:B:144:GLU:OE1	5:B:403:HOH:O	2.17	0.58
1:B:140:PHE:HE1	1:B:156:LYS:HD2	1.69	0.58
1:B:128:ILE:HD13	1:B:192:LEU:HD22	1.86	0.56
1:B:169:LYS:HD2	3:D:225:DT:H5''	1.87	0.56
1:B:52:TYR:HB2	1:B:60:VAL:HG22	1.90	0.53
1:B:33:ASN:ND2	1:B:35:ASP:H	2.06	0.53
1:B:126:PHE:CD1	1:B:196:ARG:HG3	2.46	0.51
1:B:55:LYS:HD2	1:B:55:LYS:HA	1.73	0.46
3:D:218:DC:N4	4:D:301:SO4:O3	2.40	0.46
1:B:127:PRO:HB2	1:B:197:LYS:HB3	1.96	0.46
1:B:140:PHE:CE1	1:B:156:LYS:HD2	2.50	0.45
1:B:59:ALA:HB2	1:B:74:PHE:CE2	2.52	0.44
2:C:213:DC:H2''	2:C:214:DA:H8	1.83	0.44
1:B:196:ARG:NE	5:B:415:HOH:O	2.47	0.43
1:B:71:ALA:HB2	1:B:97:TYR:CE2	2.54	0.43
1:B:144:GLU:HG2	1:B:144:GLU:O	2.18	0.42
1:B:51:GLU:HG3	1:B:61:ILE:HB	2.03	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:436:HOH:O	5:B:437:HOH:O[1_565]	2.14	0.06

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	B	185/219 (84%)	178 (96%)	7 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	B	159/188 (85%)	153 (96%)	6 (4%)	33	51

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	15	SER
1	B	33	ASN
1	B	131	GLU
1	B	144	GLU
1	B	146	GLU
1	B	169	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	33	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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	SO4	D	301	-	4,4,4	0.13	0	6,6,6	0.07	0
4	SO4	B	301	-	4,4,4	0.14	0	6,6,6	0.07	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	301	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.