



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 26, 2021 – 09:01 AM GMT

PDB ID : 1QNE  
Title : Crystal structure of the Adenovirus major late promoter TATA box bound to wild-type TBP (Arabidopsis thaliana TBP isoform 2).  
Authors : Kim, J.L.; Burley, S.K.  
Deposited on : 1999-10-14  
Resolution : 1.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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  
Xtrriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.16

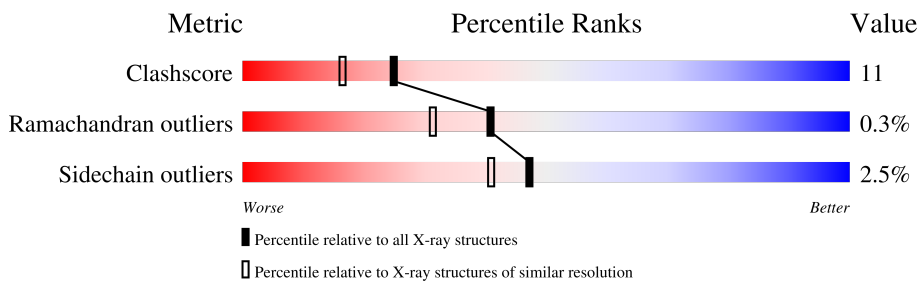
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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  $\geq 3$ , 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  $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	200	
1	B	200	
2	C	14	
2	E	14	
3	D	14	
3	F	14	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4616 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 TRANSCRIPTION INITIATION FACTOR TFIID-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	187	Total 1473	C 960	N 250	O 255	S 8	0	0	0
1	B	188	Total 1489	C 970	N 255	O 256	S 8	0	1	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	14	Total 287	C 137	N 60	O 77	P 13	0	0	0
2	E	14	Total 289	C 138	N 60	O 78	P 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	Total 279	C 136	N 44	O 86	P 13	0	0	0
3	F	14	Total 279	C 136	N 44	O 86	P 13	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	174	Total 174	O 174	0	0
4	B	180	Total 180	O 180	0	0
4	C	42	Total 42	O 42	0	0

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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
4	D	39	Total O 39 39	0	0
4	E	38	Total O 38 38	0	0
4	F	47	Total O 47 47	0	0

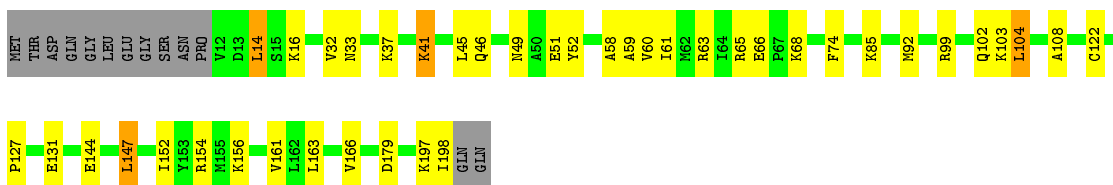
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.


Note EDS was not executed.

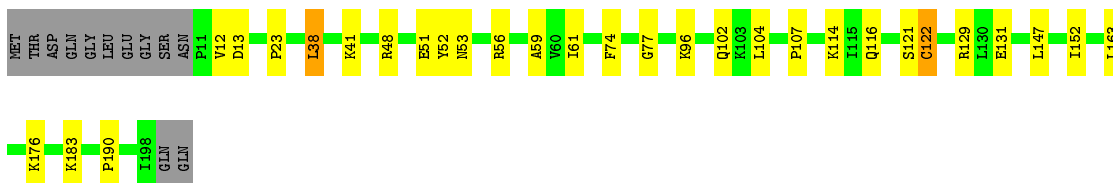
- Molecule 1: TRANSCRIPTION INITIATION FACTOR TFIIID-1

Chain A:  73% 19% 7%



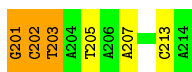
- Molecule 1: TRANSCRIPTION INITIATION FACTOR TFIIID-1

Chain B:  79% 14% 6%

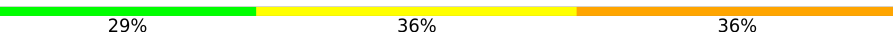


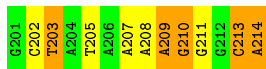
- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*GP\*GP\*CP\*A)-3')

Chain C:  57% 21% 21%



- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*GP\*GP\*CP\*A)-3')

Chain E:  29% 36% 36%



- Molecule 3: DNA (5'-D(\*TP\*GP\*CP\*CP\*CP\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain D:  14% 64% 21%



- Molecule 3: DNA (5'-D(\*TP\*GP\*CP\*CP\*CP\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain F: 21% 29% 50%



## 4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	42.00 Å 147.00 Å 57.00 Å 90.00° 90.50° 90.00°	Depositor
Resolution (Å)	6.00 – 1.90	Depositor
% Data completeness (in resolution range)	94.4 (6.00-1.90)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.197 , 0.265	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	4616	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

## 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	0.68	1/1503 (0.1%)	0.78	0/2024
1	B	0.69	1/1525 (0.1%)	0.81	0/2052
2	C	1.37	1/324 (0.3%)	1.71	7/499 (1.4%)
2	E	1.39	0/326	1.84	15/502 (3.0%)
3	D	1.39	2/310 (0.6%)	2.04	15/476 (3.2%)
3	F	1.50	6/310 (1.9%)	2.50	22/476 (4.6%)
All	All	0.96	11/4298 (0.3%)	1.33	59/6029 (1.0%)

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	CYS	CB-SG	-8.12	1.68	1.82
3	D	215	DT	C5-C7	6.27	1.53	1.50
2	C	202	DC	P-O5'	6.24	1.66	1.59
3	F	218	DC	P-O5'	5.67	1.65	1.59
3	F	222	DT	C5-C7	5.62	1.53	1.50

The worst 5 of 59 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	217	DC	O4'-C4'-C3'	-15.43	96.74	106.00
3	D	227	DG	O4'-C1'-N9	13.61	117.53	108.00
3	F	217	DC	O4'-C1'-N1	12.10	116.47	108.00
3	F	219	DC	O4'-C4'-C3'	-10.19	99.88	106.00
3	F	220	DT	C6-C5-C7	-10.15	116.81	122.90

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	1473	0	1549	36	2
1	B	1489	0	1575	27	0
2	C	287	0	154	6	0
2	E	289	0	158	4	2
3	D	279	0	162	7	0
3	F	279	0	162	11	0
4	A	174	0	0	12	1
4	B	180	0	0	10	1
4	C	42	0	0	1	0
4	D	39	0	0	1	0
4	E	38	0	0	2	0
4	F	47	0	0	0	0
All	All	4616	0	3760	84	3

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 11.

The worst 5 of 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:A:156:LYS:HA	1:A:156:LYS:HE2	1.39	1.04
1:B:41:LYS:HD3	1:B:52:TYR:HE2	1.26	1.00
1:A:66:GLU:HG2	4:A:2071:HOH:O	1.72	0.90
3:F:217:DC:H2'	3:F:218:DC:C6	2.11	0.85
1:A:198:ILE:HD11	1:B:190:PRO:HG2	1.60	0.83

All (3) 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)
4:A:2040:HOH:O	4:B:2151:HOH:O[2_647]	1.97	0.23
1:A:49:ASN:ND2	2:E:208:DA:OP1[2_647]	2.05	0.15
1:A:65:ARG:NH2	2:E:211:DG:N7[2_647]	2.11	0.09

## 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	185/200 (92%)	176 (95%)	9 (5%)	0	100	100
1	B	187/200 (94%)	179 (96%)	7 (4%)	1 (0%)	29	18
All	All	372/400 (93%)	355 (95%)	16 (4%)	1 (0%)	41	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	48	ARG

### 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	159/171 (93%)	154 (97%)	5 (3%)	40	32
1	B	162/171 (95%)	159 (98%)	3 (2%)	57	53
All	All	321/342 (94%)	313 (98%)	8 (2%)	47	41

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	LEU
1	B	104	LEU
1	B	38	LEU
1	A	104	LEU
1	A	179	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	B	53	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](#)

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.