



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:57 AM UTC

PDB ID : 1QN3 / pdb_00001qn3
Title : Crystal structure of the C(-25) Adenovirus major late promoter TATA box variant bound to wild-type TBP (Arabidopsis thaliana TBP isoform 2). TATA element recognition by the TATA box-binding protein has been conserved throughout evolution.
Authors : Patikoglou, G.A.; Kim, J.L.; Sun, L.; Yang, S.-H.; Kodadek, T.; Burley, S.K.
Deposited on : 1999-10-14
Resolution : 1.95 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 : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

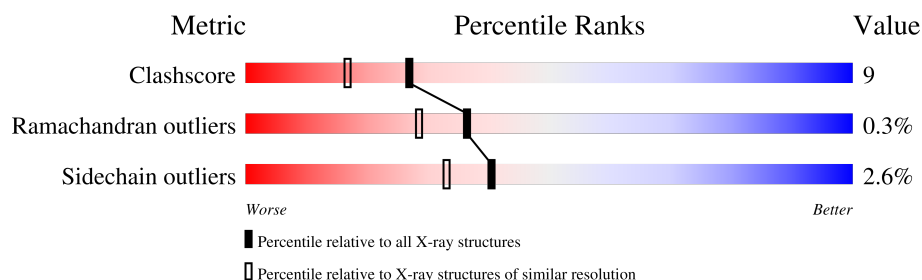
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	200	70% 19% 8%
1	B	200	72% 22% 6%
2	C	14	71% 29%
2	E	14	64% 36%
3	D	14	57% 43%
3	F	14	64% 36%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4514 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
1	A	183	Total	C	N	O	S	0	0	0
			1441	941	244	248	8			
1	B	187	Total	C	N	O	S	0	1	0
			1467	958	246	254	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	14	Total	C	N	O	P	0	0	0
			285	136	58	78	13			
2	E	14	Total	C	N	O	P	0	0	0
			287	137	58	79	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	14	Total	C	N	O	P	0	0	0
			281	136	47	85	13			
3	F	14	Total	C	N	O	P	0	0	0
			281	136	47	85	13			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	148	Total	O	0	0
			148	148		
4	B	167	Total	O	0	0
			167	167		
4	C	36	Total	O	0	0
			36	36		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	45	Total 45	O 45	0	0
4	E	36	Total 36	O 36	0	0
4	F	40	Total 40	O 40	0	0

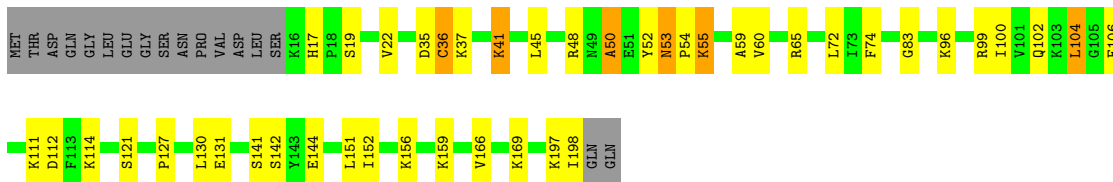
3 Residue-property plots

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.

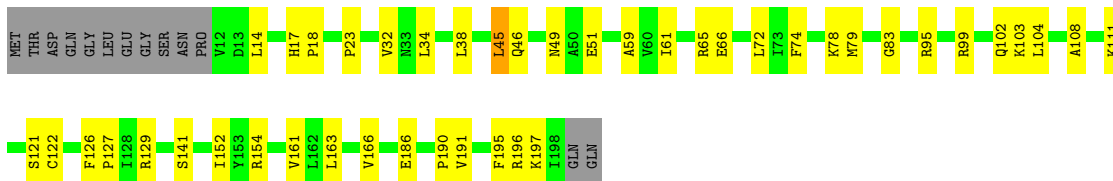
• Molecule 1: TRANSCRIPTION INITIATION FACTOR TFIIID-1

Chain A: 



• Molecule 1: TRANSCRIPTION INITIATION FACTOR TFIIID-1

Chain B: 



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

Chain C: 



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

Chain E: 



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

Chain D: 



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

Chain F: 64% 36%



4 Data and refinement statistics

Xtriage (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, α , β , γ	41.80 Å 146.70 Å 57.40 Å 90.00° 90.50° 90.00°	Depositor
Resolution (Å)	6.00 – 1.95	Depositor
% Data completeness (in resolution range)	99.5 (6.00-1.95)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.199 , 0.255	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4514	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

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	0.55	0/1471	1.04	8/1980 (0.4%)
1	B	0.54	0/1501	0.97	6/2022 (0.3%)
2	C	0.56	0/321	1.00	0/494
2	E	0.51	0/323	1.03	0/497
3	D	0.59	0/313	1.33	1/481 (0.2%)
3	F	0.65	0/313	1.31	1/481 (0.2%)
All	All	0.55	0/4242	1.06	16/5955 (0.3%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	ASN	CA-C-N	8.42	128.92	119.32
1	A	53	ASN	C-N-CA	8.42	128.92	119.32
1	A	50	ALA	N-CA-C	8.12	121.52	110.55
1	B	111	LYS	N-CA-C	6.45	120.17	109.46
3	D	220	DG	C4'-C3'-O3'	-6.31	100.53	110.00
1	B	83	GLY	N-CA-C	6.10	121.95	114.69
1	B	141	SER	N-CA-C	5.75	119.33	110.42
1	B	191	VAL	N-CA-C	-5.61	105.38	110.82
3	F	216	DG	P-O3'-C3'	5.53	128.49	120.20
1	A	83	GLY	N-CA-C	5.48	122.34	115.21
1	A	48	ARG	N-CA-C	-5.42	106.67	113.28
1	A	141	SER	N-CA-C	5.37	118.56	109.76
1	B	17	HIS	CA-C-N	5.22	125.27	119.32
1	B	17	HIS	C-N-CA	5.22	125.27	119.32
1	A	130	LEU	N-CA-C	5.11	116.53	111.07
1	A	72	LEU	N-CA-C	-5.04	100.19	108.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1441	0	1518	27	1
1	B	1467	0	1537	25	3
2	C	285	0	154	4	2
2	E	287	0	158	7	0
3	D	281	0	161	7	0
3	F	281	0	161	5	0
4	A	148	0	0	7	1
4	B	167	0	0	9	1
4	C	36	0	0	2	0
4	D	45	0	0	2	0
4	E	36	0	0	3	0
4	F	40	0	0	2	0
All	All	4514	0	3689	71	4

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 9.

All (71) 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:66:GLU:HG2	4:B:2052:HOH:O	1.57	1.05
1:B:46:GLN:NE2	4:B:2025:HOH:O	1.95	0.99
1:A:144:GLU:HG2	4:A:2106:HOH:O	1.65	0.95
3:D:215:DT:H2''	3:D:216:DG:H5'	1.48	0.93
2:E:201:DG:C4	4:E:2007:HOH:O	2.31	0.83
1:B:65:ARG:HD3	4:B:2049:HOH:O	1.78	0.83
1:A:52:TYR:CZ	1:A:54:PRO:HG3	2.15	0.81
1:B:66:GLU:CG	4:B:2052:HOH:O	2.21	0.81
1:A:65:ARG:NH1	4:A:2047:HOH:O	2.23	0.71
1:A:198:ILE:HD11	1:B:190:PRO:HG2	1.74	0.70
3:F:215:DT:O5'	4:F:2009:HOH:O	2.10	0.68
3:D:215:DT:C2	4:D:2008:HOH:O	2.47	0.67
3:D:215:DT:H2'	3:D:216:DG:C8	2.29	0.67
1:B:127:PRO:HB2	1:B:197:LYS:HB3	1.82	0.62
1:B:59:ALA:HB2	1:B:74:PHE:CE1	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:201:DG:N3	4:E:2007:HOH:O	2.33	0.59
1:A:96:LYS:O	1:A:100:ILE:HG13	2.02	0.59
1:B:102:GLN:HG3	1:B:108:ALA:HB3	1.84	0.58
1:A:111:LYS:HG2	1:A:112:ASP:OD2	2.03	0.58
3:D:215:DT:H4'	3:D:216:DG:OP1	2.03	0.58
2:C:201:DG:H2'	2:C:202:DC:C6	2.39	0.58
1:A:41:LYS:HE2	1:A:41:LYS:HA	1.86	0.57
1:B:152:ILE:HD12	1:B:163:LEU:HD22	1.86	0.57
1:A:19:SER:HA	1:B:14:LEU:HB2	1.88	0.56
1:A:53:ASN:HB3	1:A:55:LYS:NZ	2.21	0.55
1:B:61:ILE:HG12	1:B:72:LEU:HD23	1.88	0.55
1:A:121:SER:HB2	4:A:2129:HOH:O	2.07	0.53
3:D:215:DT:N3	4:D:2008:HOH:O	2.42	0.52
1:B:34:LEU:HD12	1:B:38:LEU:HD11	1.90	0.52
1:A:52:TYR:CE2	1:A:54:PRO:HG3	2.45	0.52
1:A:111:LYS:HG2	1:A:112:ASP:CG	2.33	0.52
1:A:131:GLU:HB3	4:A:2096:HOH:O	2.10	0.52
2:E:210:DG:H2''	2:E:211:DG:C8	2.46	0.51
1:B:126:PHE:CD1	1:B:196:ARG:HG3	2.45	0.51
1:A:102:GLN:HA	1:A:106:PHE:O	2.11	0.50
1:B:129:ARG:NH2	1:B:195:PHE:HA	2.27	0.49
1:A:50:ALA:HB1	1:A:60:VAL:HG13	1.94	0.49
1:B:99:ARG:HG3	1:B:103:LYS:HE3	1.93	0.49
2:E:202:DC:H2'	2:E:203:DT:C6	2.48	0.49
1:A:142:SER:HB2	1:A:152:ILE:HB	1.95	0.48
2:C:210:DG:H2''	4:C:2033:HOH:O	2.12	0.48
1:A:127:PRO:HA	1:A:166:VAL:O	2.12	0.48
3:D:217:DC:H2'	3:D:218:DC:C6	2.49	0.47
1:A:17:HIS:CE1	1:A:22:VAL:HG13	2.50	0.46
1:B:49:ASN:N	4:B:2028:HOH:O	2.41	0.46
3:F:219:DC:OP2	3:F:219:DC:H6	1.99	0.45
1:A:156:LYS:HD3	4:A:2117:HOH:O	2.16	0.45
4:A:2062:HOH:O	2:C:210:DG:H5''	2.15	0.45
1:B:154:ARG:HG2	1:B:161:VAL:HG22	1.98	0.44
1:B:127:PRO:HA	1:B:166:VAL:O	2.17	0.44
1:A:104:LEU:HD12	1:A:104:LEU:HA	1.85	0.44
4:B:2142:HOH:O	3:F:225:DT:H5'	2.17	0.44
2:C:210:DG:C2'	4:C:2033:HOH:O	2.64	0.44
3:F:227:DG:H5'	4:F:2039:HOH:O	2.18	0.44
1:A:197:LYS:HE3	4:A:2146:HOH:O	2.18	0.43
1:B:126:PHE:HB2	1:B:196:ARG:NE	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:PRO:HD2	1:B:186:GLU:CD	2.43	0.43
2:E:210:DG:N7	3:F:219:DC:N3	2.67	0.42
1:A:41:LYS:O	1:A:45:LEU:HG	2.19	0.42
1:A:99:ARG:HD3	1:A:102:GLN:OE1	2.19	0.42
1:B:32:VAL:CG1	1:B:79:MET:HE2	2.50	0.42
2:E:203:DT:H5'	4:E:2008:HOH:O	2.19	0.42
1:A:151:LEU:C	1:A:151:LEU:HD23	2.46	0.41
1:B:121:SER:HB2	4:B:2142:HOH:O	2.21	0.41
1:A:59:ALA:HB2	1:A:74:PHE:CE2	2.56	0.41
2:E:210:DG:H2'	2:E:210:DG:N3	2.35	0.41
1:A:169:LYS:HG2	3:D:225:DT:H5''	2.03	0.41
1:B:23:PRO:HA	1:B:122:CYS:HB3	2.01	0.41
1:B:78:LYS:NZ	4:B:2069:HOH:O	2.54	0.41
1:B:95:ARG:NH1	4:B:2082:HOH:O	2.53	0.41
1:A:35:ASP:O	1:A:36:CYS:HB3	2.20	0.40

All (4) 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:2126:HOH:O	4:B:2024:HOH:O[2_646]	1.56	0.64
1:B:65:ARG:NH1	2:C:211:DG:O6[2_656]	1.81	0.39
1:B:65:ARG:NH2	2:C:211:DG:N7[2_656]	2.00	0.20
1:A:159:LYS:CD	1:B:45:LEU:O[2_646]	2.09	0.11

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	181/200 (90%)	177 (98%)	3 (2%)	1 (1%)	21	12
1	B	186/200 (93%)	182 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	367/400 (92%)	359 (98%)	7 (2%)	1 (0%)	36 28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	CYS

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	155/171 (91%)	150 (97%)	5 (3%)	34 25
1	B	158/171 (92%)	155 (98%)	3 (2%)	50 45
All	All	313/342 (92%)	305 (97%)	8 (3%)	40 33

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LYS
1	A	41	LYS
1	A	55	LYS
1	A	104	LEU
1	A	114	LYS
1	B	45	LEU
1	B	51	GLU
1	B	104	LEU

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	26	GLN
1	B	46	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 oligosaccharides 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.