



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:01 PM UTC

PDB ID : 3DOM / pdb_00003dom
Title : Crystal Structure of the complex between Tfb5 and the C-terminal domain of Tfb2
Authors : Kainov, D.E.; Cavarelli, J.; Egly, J.M.; Poterszman, A.
Deposited on : 2008-07-04
Resolution : 2.60 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

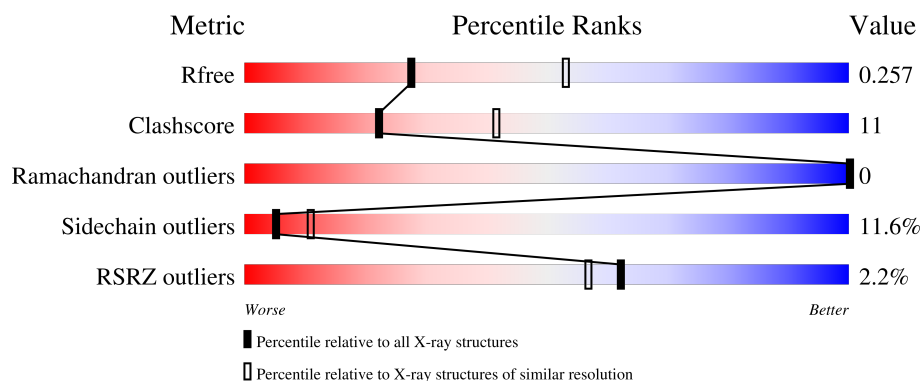
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 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 $\leq 5\%$. 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	108	% 46% 19% • 32%
1	C	108	49% 10% 9% • 29%
2	B	71	6% 66% 13% • 18%
2	D	71	% 69% 13% 10% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2323 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 RNA polymerase II transcription factor B subunit 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	73	Total	C	N	O	0	0	0
			615	399	100	116			
1	C	77	Total	C	N	O	0	0	0
			643	417	104	122			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	406	GLY	-	expression tag	UNP Q02939
A	407	PRO	-	expression tag	UNP Q02939
A	408	HIS	-	expression tag	UNP Q02939
A	409	MET	-	expression tag	UNP Q02939
A	410	ALA	-	expression tag	UNP Q02939
A	411	SER	-	expression tag	UNP Q02939
C	406	GLY	-	expression tag	UNP Q02939
C	407	PRO	-	expression tag	UNP Q02939
C	408	HIS	-	expression tag	UNP Q02939
C	409	MET	-	expression tag	UNP Q02939
C	410	ALA	-	expression tag	UNP Q02939
C	411	SER	-	expression tag	UNP Q02939

- Molecule 2 is a protein called RNA polymerase II transcription factor B subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	58	Total	C	N	O	S	0	0	0
			454	287	80	85	2			
2	D	65	Total	C	N	O	S	0	0	0
			514	326	90	95	3			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	32	Total 32	O 32	0	0
3	B	20	Total 20	O 20	0	0
3	C	30	Total 30	O 30	0	0
3	D	15	Total 15	O 15	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	37.58Å 103.59Å 114.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.70 – 2.60 30.70 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.0 (30.70-2.60) 95.9 (30.70-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.71 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.202 , 0.261 0.203 , 0.257	Depositor DCC
R_{free} test set	703 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 28.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2323	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.21% 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

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	0.97	0/626	1.17	3/840 (0.4%)
1	C	1.05	2/656 (0.3%)	1.38	9/885 (1.0%)
2	B	0.93	0/458	1.14	0/617
2	D	0.84	1/520 (0.2%)	1.06	0/701
All	All	0.96	3/2260 (0.1%)	1.21	12/3043 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	451	VAL	CA-CB	5.53	1.60	1.54
1	C	464	THR	N-CA	-5.40	1.39	1.46
2	D	13	ASP	CA-C	5.37	1.59	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	435	PRO	N-CA-C	-14.19	90.62	114.75
1	A	439	ASP	N-CA-C	-11.60	99.50	112.72
1	C	434	PRO	CB-CA-C	-7.21	102.13	110.92
1	A	440	GLN	N-CA-C	-6.07	104.75	111.36
1	A	488	LYS	N-CA-C	-5.70	106.31	113.72
1	C	434	PRO	CA-C-N	5.68	126.55	120.47
1	C	434	PRO	C-N-CA	5.68	126.55	120.47
1	C	433	LEU	CA-C-N	-5.60	114.61	120.38
1	C	433	LEU	C-N-CA	-5.60	114.61	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	488	LYS	N-CA-C	-5.32	106.84	113.38
1	C	466	GLN	N-CA-C	5.27	117.45	111.02
1	C	498	ASN	N-CA-C	5.22	117.39	111.02

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	434	PRO	Peptide
1	C	463	GLU	Peptide

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	615	0	619	15	0
1	C	643	0	644	18	0
2	B	454	0	480	10	0
2	D	514	0	541	6	0
3	A	32	0	0	2	0
3	B	20	0	0	2	0
3	C	30	0	0	0	0
3	D	15	0	0	0	0
All	All	2323	0	2284	48	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 11.

All (48) 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:485:ASP:OD2	1:C:488:LYS:HD3	1.51	1.07
1:C:464:THR:HG22	1:C:467:GLU:H	1.35	0.91
1:C:463:GLU:HB2	1:C:467:GLU:OE2	1.79	0.82
1:A:485:ASP:OD2	1:A:488:LYS:HD2	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:38:THR:HG22	2:D:39:HIS:CE1	2.23	0.73
1:A:438:VAL:HG12	1:A:441:ILE:HG13	1.74	0.69
1:C:471:LEU:HD12	1:C:501:VAL:HG13	1.76	0.67
1:C:448:LEU:O	1:C:451:VAL:HG13	1.97	0.65
1:C:435:PRO:CD	1:C:436:THR:H	1.99	0.64
1:C:442:ARG:O	1:C:446:LEU:HD22	1.98	0.64
1:C:446:LEU:O	1:C:450:ARG:HG3	2.00	0.60
1:A:489:LYS:HE3	3:A:61:HOH:O	2.02	0.58
2:B:5:ARG:HD2	3:B:75:HOH:O	2.04	0.56
1:A:438:VAL:CG1	1:A:441:ILE:HG13	2.37	0.55
1:C:435:PRO:HD2	1:C:436:THR:H	1.69	0.55
1:C:464:THR:CG2	1:C:467:GLU:H	2.14	0.53
1:C:471:LEU:CD1	1:C:501:VAL:HG13	2.38	0.53
1:C:503:ASP:OD1	1:C:507:ARG:HD2	2.09	0.52
2:B:27:MET:O	2:B:28:SER:OG	2.25	0.52
1:A:437:VAL:HG13	1:A:438:VAL:N	2.24	0.51
2:B:24:ASP:O	2:B:28:SER:N	2.42	0.51
1:C:464:THR:HG22	1:C:467:GLU:N	2.16	0.51
2:D:27:MET:HB3	2:D:29:ASP:OD1	2.10	0.51
1:A:491:PHE:HE2	1:A:493:ILE:HD12	1.77	0.49
1:A:454:TYR:CG	1:A:482:LEU:HD13	2.48	0.49
1:C:435:PRO:CD	1:C:436:THR:N	2.67	0.48
1:A:493:ILE:HD13	1:A:501:VAL:HG11	1.96	0.47
2:D:19:LEU:O	2:D:22:GLN:HB2	2.14	0.47
1:A:438:VAL:C	1:A:440:GLN:H	2.19	0.47
1:A:507:ARG:HG2	1:A:507:ARG:HH11	1.81	0.46
1:C:495:LYS:HB3	1:C:495:LYS:HE3	1.53	0.46
1:C:493:ILE:HD11	1:C:498:ASN:HA	1.97	0.45
1:C:471:LEU:HD12	1:C:501:VAL:CG1	2.46	0.45
1:A:458:LEU:HD22	2:B:41:LEU:HD22	1.98	0.44
2:B:36:ASP:C	2:B:36:ASP:OD1	2.61	0.44
2:B:48:GLU:OE1	2:B:48:GLU:HA	2.18	0.43
2:D:35:LEU:HA	2:D:35:LEU:HD23	1.73	0.42
2:D:60:LYS:HA	2:D:60:LYS:HD3	1.78	0.42
3:A:11:HOH:O	2:B:2:ALA:HB2	2.19	0.42
1:C:464:THR:CG2	1:C:466:GLN:HB2	2.50	0.42
1:A:438:VAL:C	1:A:440:GLN:N	2.76	0.42
2:B:58:LEU:O	2:B:59:SER:OG	2.29	0.42
2:B:34:GLU:CD	3:B:86:HOH:O	2.62	0.42
2:D:53:GLU:OE1	2:D:56:ARG:NH1	2.53	0.42
1:A:470:LEU:HD23	1:A:509:LEU:HD21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:475:ALA:HB2	1:A:493:ILE:HD11	2.03	0.41
2:B:24:ASP:O	2:B:28:SER:HA	2.21	0.41
1:A:456:GLY:HA3	1:A:493:ILE:O	2.21	0.41

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	71/108 (66%)	70 (99%)	1 (1%)	0	100	100
1	C	75/108 (69%)	73 (97%)	2 (3%)	0	100	100
2	B	56/71 (79%)	53 (95%)	3 (5%)	0	100	100
2	D	63/71 (89%)	55 (87%)	8 (13%)	0	100	100
All	All	265/358 (74%)	251 (95%)	14 (5%)	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	A	68/100 (68%)	63 (93%)	5 (7%)	13	29
1	C	72/100 (72%)	62 (86%)	10 (14%)	3	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	52/65 (80%)	50 (96%)	2 (4%)	29	56
2	D	59/65 (91%)	47 (80%)	12 (20%)	1	2
All	All	251/330 (76%)	222 (88%)	29 (12%)	5	11

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	463	GLU
1	A	481	LEU
1	A	487	LYS
1	A	496	GLU
1	A	501	VAL
2	B	27	MET
2	B	58	LEU
1	C	436	THR
1	C	439	ASP
1	C	442	ARG
1	C	446	LEU
1	C	451	VAL
1	C	464	THR
1	C	467	GLU
1	C	495	LYS
1	C	498	ASN
1	C	502	LEU
2	D	3	ARG
2	D	13	ASP
2	D	19	LEU
2	D	22	GLN
2	D	26	LYS
2	D	27	MET
2	D	28	SER
2	D	38	THR
2	D	55	ASN
2	D	56	ARG
2	D	58	LEU
2	D	60	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	22	GLN
1	C	445	GLN
1	C	498	ASN
2	D	11	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 [i](#)

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

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, 95th 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 ²)	Q<0.9
1	A	73/108 (67%)	-0.24	1 (1%) 73 69	2, 10, 18, 24	0
1	C	77/108 (71%)	-0.24	0 100 100	2, 10, 19, 23	0
2	B	58/71 (81%)	0.61	4 (6%) 23 18	5, 10, 17, 21	0
2	D	65/71 (91%)	0.10	1 (1%) 72 68	4, 11, 20, 27	0
All	All	273/358 (76%)	0.02	6 (2%) 62 57	2, 10, 19, 27	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	2	ALA	3.4
2	D	26	LYS	2.8
1	A	461	ASP	2.7
2	B	36	ASP	2.4
2	B	49	PHE	2.0
2	B	52	HIS	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 oligosaccharides 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.