



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 04:47 AM UTC

PDB ID : 5NV6 / pdb_00005nv6
Title : Structure of human transforming growth factor beta-induced protein (TGF- β 1p).
Authors : Garcia-Castellanos, R.; Nielsen, S.N.; Runager, K.; Thogersen, B.I.; Goulas, T.; Enghild, J.J.; Gomis-Ruth, F.X.
Deposited on : 2017-05-03
Resolution : 2.93 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

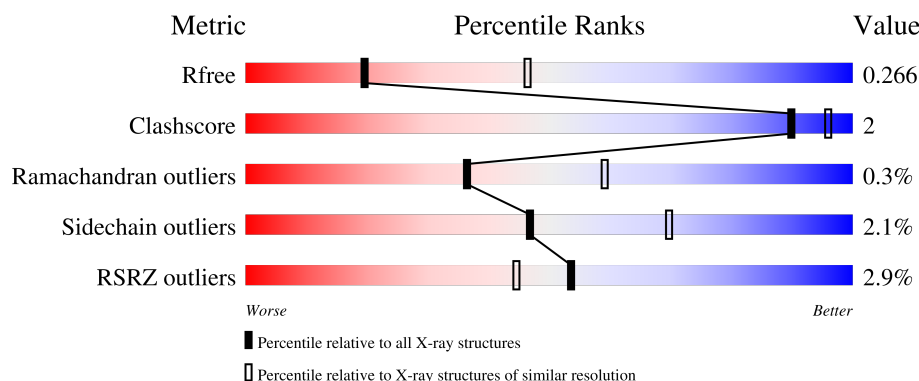
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.93 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	1159 (2.96-2.92)
Clashscore	190562	1184 (2.96-2.92)
Ramachandran outliers	187476	1131 (2.96-2.92)
Sidechain outliers	187428	1131 (2.96-2.92)
RSRZ outliers	180081	1159 (2.96-2.92)

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	683	2% 80% 7% 13%
1	B	683	3% 79% 7% 14%

2 Entry composition [i](#)

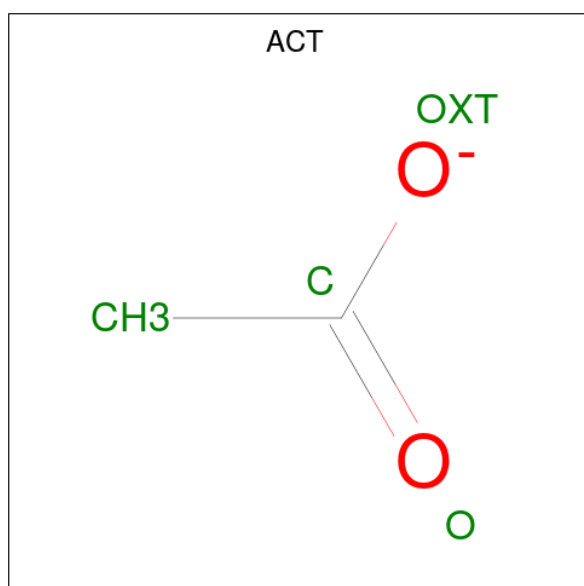
There are 3 unique types of molecules in this entry. The entry contains 9273 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Transforming growth factor-beta-induced protein ig-h3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	S	0	0	0
			4559	2866	797	873	23			
1	B	589	Total	C	N	O	S	0	0	0
			4516	2841	787	865	23			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	79	Total	O	0	0
			79	79		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	115	Total	O	0	0
			115	115		

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	114.84Å 114.84Å 181.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.93 25.00 – 2.93	Depositor EDS
% Data completeness (in resolution range)	98.5 (25.00-2.93) 98.3 (25.00-2.93)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.91Å)	Xtriage
Refinement program	PHENIX, BUSTER-TNT	Depositor
R, R_{free}	0.219 , 0.263 0.223 , 0.266	Depositor DCC
R_{free} test set	652 reflections (2.24%)	wwPDB-VP
Wilson B-factor (Å ²)	61.4	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.047 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9273	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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

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

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	0.78	0/4635	1.28	15/6296 (0.2%)
1	B	0.79	1/4590 (0.0%)	1.28	13/6234 (0.2%)
All	All	0.79	1/9225 (0.0%)	1.28	28/12530 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	500	PRO	CA-C	5.42	1.54	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	362	GLY	N-CA-C	6.39	119.49	110.74
1	A	232	ILE	CA-C-N	6.38	129.46	120.28
1	A	232	ILE	C-N-CA	6.38	129.46	120.28
1	B	232	ILE	CA-C-N	6.35	129.43	120.28
1	B	232	ILE	C-N-CA	6.35	129.43	120.28
1	B	362	GLY	N-CA-C	6.35	119.43	110.74
1	B	483	ASP	CA-CB-CG	5.95	118.55	112.60
1	B	513	ASN	CA-CB-CG	5.83	118.43	112.60
1	B	227	GLY	N-CA-C	5.82	118.01	110.45
1	A	227	GLY	N-CA-C	5.72	117.89	110.45
1	A	270	GLU	CA-C-N	5.69	125.50	120.10
1	A	270	GLU	C-N-CA	5.69	125.50	120.10
1	A	450	TYR	CA-C-N	5.63	128.15	120.54
1	A	450	TYR	C-N-CA	5.63	128.15	120.54
1	A	272	ASN	CA-CB-CG	5.51	118.11	112.60
1	B	270	GLU	CA-C-N	5.47	125.30	120.10
1	B	270	GLU	C-N-CA	5.47	125.30	120.10
1	B	450	TYR	CA-C-N	5.30	129.62	121.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	450	TYR	C-N-CA	5.30	129.62	121.19
1	B	561	ASP	CA-C-N	5.20	128.21	120.31
1	B	561	ASP	C-N-CA	5.20	128.21	120.31
1	A	125	THR	CA-C-N	5.16	129.66	122.08
1	A	125	THR	C-N-CA	5.16	129.66	122.08
1	A	73	ILE	CA-C-N	5.04	131.17	121.54
1	A	73	ILE	C-N-CA	5.04	131.17	121.54
1	A	400	LEU	CA-C-N	5.01	125.64	120.03
1	A	400	LEU	C-N-CA	5.01	125.64	120.03
1	B	492	PHE	CA-CB-CG	5.00	118.81	113.80

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	4559	0	4611	19	0
1	B	4516	0	4575	14	0
2	A	4	0	3	0	0
3	A	79	0	0	0	0
3	B	115	0	0	0	0
All	All	9273	0	9189	33	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 2.

All (33) 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:110:LEU:HD11	1:A:132:MET:HE1	1.82	0.60
1:A:128:LEU:HD12	1:A:168:LEU:HD12	1.83	0.60
1:A:122:THR:HG23	1:A:128:LEU:HB3	1.85	0.58
1:B:221:ASP:H	1:B:230:HIS:CE1	2.22	0.58
1:A:217:LEU:HD11	1:A:230:HIS:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LEU:HD11	1:B:132:MET:HE1	1.90	0.53
1:B:398:ALA:HB1	1:B:432:ARG:HG3	1.89	0.53
1:A:476:ASN:HB3	1:A:619:MET:HE3	1.90	0.52
1:B:449:LEU:HA	1:B:453:GLN:OE1	2.10	0.51
1:A:618:ILE:HB	1:A:625:VAL:HB	1.96	0.47
1:A:276:THR:HB	1:A:312:LEU:HB2	1.98	0.46
1:B:618:ILE:HB	1:B:625:VAL:HB	1.96	0.46
1:A:379:LEU:HB3	1:A:404:LEU:HA	1.97	0.46
1:A:579:VAL:HG12	1:A:619:MET:HE1	1.96	0.46
1:B:256:LEU:HD13	1:B:280:PRO:HD2	1.98	0.46
1:B:276:THR:HB	1:B:312:LEU:HB2	1.98	0.46
1:B:506:MET:HE2	1:B:528:THR:HG23	1.99	0.44
1:B:248:GLU:HA	1:B:257:ARG:HD3	2.00	0.44
1:A:484:LYS:HB2	1:A:491:LEU:HB3	2.00	0.43
1:A:553:ARG:O	1:A:557:ARG:HB2	2.18	0.43
1:B:285:PHE:HD1	1:B:293:LEU:HD11	1.83	0.43
1:A:451:HIS:HA	1:A:467:VAL:HG12	2.00	0.43
1:A:256:LEU:HD13	1:A:280:PRO:HD2	1.99	0.43
1:A:248:GLU:HA	1:A:257:ARG:HD3	2.01	0.43
1:B:476:ASN:HB3	1:B:619:MET:HE3	2.01	0.42
1:A:477:SER:OG	1:A:494:MET:HB3	2.20	0.42
1:A:379:LEU:HD21	1:A:412:LEU:HB2	2.02	0.42
1:A:468:TYR:HB2	1:A:471:SER:HB2	2.01	0.42
1:A:539:VAL:HG22	1:A:573:ILE:HG23	2.01	0.41
1:B:468:TYR:HB2	1:B:471:SER:HB2	2.03	0.41
1:A:276:THR:OG1	1:A:315:ALA:HA	2.22	0.40
1:B:379:LEU:HD21	1:B:412:LEU:HB2	2.03	0.40
1:B:122:THR:HG23	1:B:128:LEU:HB3	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	593/683 (87%)	567 (96%)	24 (4%)	2 (0%)	36 59
1	B	587/683 (86%)	566 (96%)	20 (3%)	1 (0%)	43 65
All	All	1180/1366 (86%)	1133 (96%)	44 (4%)	3 (0%)	36 59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	70	GLN
1	A	74	CYS
1	A	460	GLY

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	505/574 (88%)	496 (98%)	9 (2%)	51 72
1	B	501/574 (87%)	489 (98%)	12 (2%)	43 67
All	All	1006/1148 (88%)	985 (98%)	21 (2%)	47 70

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	SER
1	A	128	LEU
1	A	172	ARG
1	A	189	MET
1	A	322	VAL
1	A	342	ASP
1	A	472	LEU
1	A	553	ARG
1	A	633	GLN
1	B	63	THR
1	B	67	GLN
1	B	68	TRP
1	B	115	SER

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Mol	Chain	Res	Type
1	B	117	THR
1	B	172	ARG
1	B	313	LYS
1	B	355	LYS
1	B	387	ASP
1	B	483	ASP
1	B	513	ASN
1	B	596	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	145	ASN
1	A	201	GLN
1	A	222	HIS
1	A	230	HIS
1	A	241	ASN
1	A	245	GLN
1	A	274	GLN
1	A	451	HIS
1	A	633	GLN
1	B	201	GLN
1	B	203	HIS
1	B	223	HIS
1	B	241	ASN
1	B	245	GLN
1	B	402	ASN
1	B	433	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

1 ligand is 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	$\# Z > 2$
2	ACT	A	701	-	3,3,3	1.22	0	3,3,3	0.92	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.

No monomer is involved in short contacts.

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, 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	595/683 (87%)	0.31	15 (2%)	58 49	42, 69, 99, 135	0
1	B	589/683 (86%)	0.38	19 (3%)	50 42	27, 64, 130, 176	0
All	All	1184/1366 (86%)	0.34	34 (2%)	53 44	27, 68, 119, 176	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	241	ASN	3.2
1	A	100	ALA	2.8
1	A	427	ILE	2.7
1	B	69	TYR	2.7
1	A	636	ALA	2.6
1	A	68	TRP	2.6
1	B	51	VAL	2.6
1	B	635	PRO	2.6
1	A	546	ALA	2.6
1	A	353	SER	2.5
1	A	46	PRO	2.5
1	B	103	LEU	2.5
1	B	151	LEU	2.5
1	A	69	TYR	2.4
1	B	547	PHE	2.4
1	B	163	VAL	2.3
1	A	543	THR	2.3
1	B	80	ILE	2.3
1	B	627	VAL	2.3
1	A	637	ASN	2.3
1	B	100	ALA	2.3
1	B	48	VAL	2.2
1	B	65	CYS	2.2
1	B	271	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	80	ILE	2.2
1	A	344	LEU	2.1
1	A	66	LYS	2.1
1	A	65	CYS	2.1
1	A	180	VAL	2.1
1	B	68	TRP	2.1
1	B	165	ILE	2.1
1	B	614	ALA	2.1
1	B	128	LEU	2.0
1	B	62	PHE	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](#)

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. The B-factors column lists the minimum, median, 95th 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	B-factors($\text{\AA}^2$)	Q<0.9
2	ACT	A	701	4/4	0.71	0.14	65,65,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.