



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:15 PM UTC

PDB ID : 1WWC / pdb_00001wwc
Title : NT3 BINDING DOMAIN OF HUMAN TRKC RECEPTOR
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Deposited on : 1999-04-30
Resolution : 1.90 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

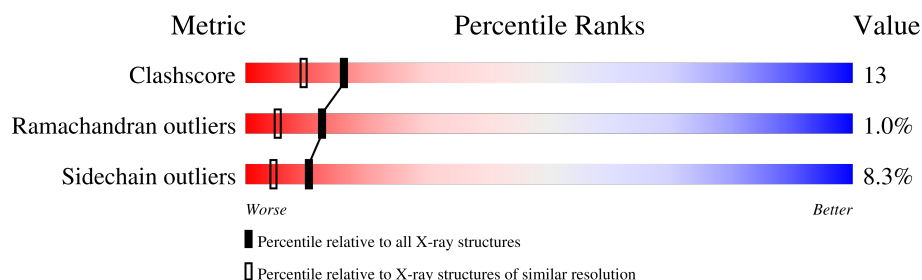
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (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 ≥ 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	118	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 958 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 PROTEIN (NT-3 GROWTH FACTOR RECEPTOR TRKC).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	105	Total	C	N	O	S	0	0	0
			857	551	146	158	2			

There are 9 discrepancies between the modelled and reference sequences:

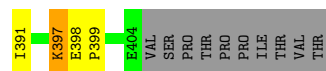
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	SEE REMARK 999	UNP Q16288
A	?	-	SER	SEE REMARK 999	UNP Q16288
A	?	-	THR	SEE REMARK 999	UNP Q16288
A	?	-	ASP	SEE REMARK 999	UNP Q16288
A	?	-	ASN	SEE REMARK 999	UNP Q16288
A	?	-	PHE	SEE REMARK 999	UNP Q16288
A	?	-	ILE	SEE REMARK 999	UNP Q16288
A	?	-	LEU	SEE REMARK 999	UNP Q16288
A	402	VAL	PHE	engineered mutation	UNP Q16288

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	101	Total	O	0	0
			101	101		

Note EDS was not executed.

- Molecule 1: PROTEIN (NT-3 GROWTH FACTOR RECEPTOR TRKC)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	82.90 Å 82.90 Å 30.40 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.90	Depositor
% Data completeness (in resolution range)	94.7 (20.00-1.90)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.187 , 0.288	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	958	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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	0.97	0/885	1.20	5/1210 (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	A	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	363	LEU	N-CA-C	-7.11	97.65	108.96
1	A	346	LYS	N-CA-C	-7.05	104.81	113.41
1	A	353	TYR	N-CA-C	-6.24	105.70	113.38
1	A	304	PRO	CA-C-N	5.44	125.38	119.78
1	A	304	PRO	C-N-CA	5.44	125.38	119.78

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	353	TYR	Sidechain

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	857	0	822	22	0
2	A	101	0	0	4	0
All	All	958	0	822	22	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 13.

All (22) 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:381:LYS:HG3	1:A:386:THR:HG22	1.18	1.12
1:A:318:GLU:OE2	1:A:366:ASN:HB2	1.94	0.67
1:A:357:GLU:HB2	2:A:468:HOH:O	1.98	0.63
1:A:314:GLU:HG2	2:A:459:HOH:O	2.00	0.61
1:A:347:ILE:HG13	1:A:348:ILE:HD12	1.81	0.61
1:A:347:ILE:HG22	2:A:434:HOH:O	2.02	0.59
1:A:370:HIS:CD2	1:A:398:GLU:OE2	2.56	0.59
1:A:397:LYS:HE2	1:A:399:PRO:HA	1.90	0.54
1:A:321:ILE:HD13	1:A:391:ILE:HD11	1.91	0.52
1:A:337:HIS:HB2	1:A:342:LEU:HD13	1.95	0.49
1:A:381:LYS:HG2	1:A:382:ASN:N	2.29	0.48
1:A:300:THR:HG22	1:A:301:VAL:N	2.28	0.48
1:A:324:VAL:HG22	1:A:360:GLU:HG3	1.96	0.48
1:A:381:LYS:HG3	1:A:386:THR:CG2	2.13	0.47
1:A:301:VAL:O	1:A:301:VAL:HG23	2.15	0.47
1:A:332:THR:O	1:A:380:ALA:HA	2.15	0.46
1:A:326:ARG:HA	1:A:357:GLU:O	2.16	0.45
1:A:378:LEU:O	1:A:388:ASN:HA	2.17	0.45
1:A:355:GLU:HG3	2:A:480:HOH:O	2.19	0.43
1:A:318:GLU:HG2	1:A:366:ASN:HA	2.01	0.41
1:A:347:ILE:O	1:A:347:ILE:HD12	2.21	0.41
1:A:330:PRO:HA	1:A:331:PRO:HD3	1.93	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	103/118 (87%)	100 (97%)	2 (2%)	1 (1%)	12 5

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	345	SER

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	96/108 (89%)	88 (92%)	8 (8%)	10 4

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

Mol	Chain	Res	Type
1	A	315	LEU
1	A	317	LEU
1	A	333	LEU
1	A	344	GLU
1	A	347	ILE
1	A	367	LYS
1	A	381	LYS
1	A	397	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	334	HIS
1	A	370	HIS
1	A	388	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 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.