



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 02:34 AM UTC

PDB ID : 4DI3 / pdb_00004di3
Title : Crystal structure of a 2:1 complex of Treponema pallidum TatP(T) (Tp0957)
bound to TatT (Tp0956)
Authors : Brautigam, C.A.; Deka, R.K.; Norgard, M.V.
Deposited on : 2012-01-30
Resolution : 3.05 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

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

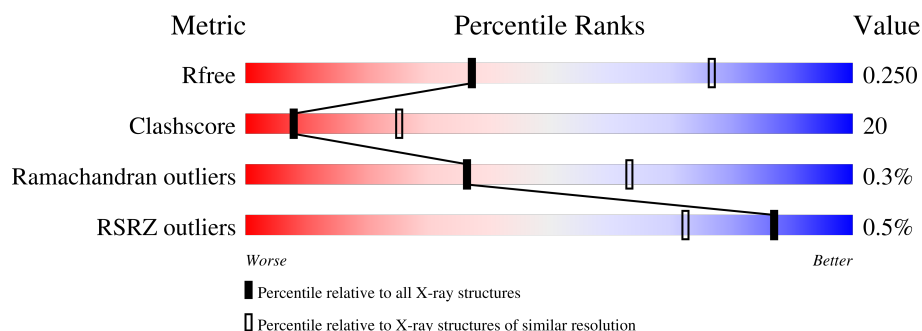
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	D	324	% 52% 40% 6%
1	E	324	% 55% 41% •
2	A	301	66% 24% 9%
2	B	301	67% 24% 9%
2	C	301	% 62% 28% • 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11376 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 TatP(T) (Tp0957).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	303	Total	C	N	O	S	0	0	0
			2385	1527	416	431	11			
1	E	317	Total	C	N	O	S	0	0	0
			2493	1594	435	453	11			

There are 6 discrepancies between the modelled and reference sequences:

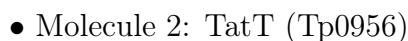
Chain	Residue	Modelled	Actual	Comment	Reference
D	5	GLY	-	expression tag	UNP O83923
D	6	ARG	-	expression tag	UNP O83923
D	185	VAL	ALA	engineered mutation	UNP O83923
E	5	GLY	-	expression tag	UNP O83923
E	6	ARG	-	expression tag	UNP O83923
E	185	VAL	ALA	engineered mutation	UNP O83923

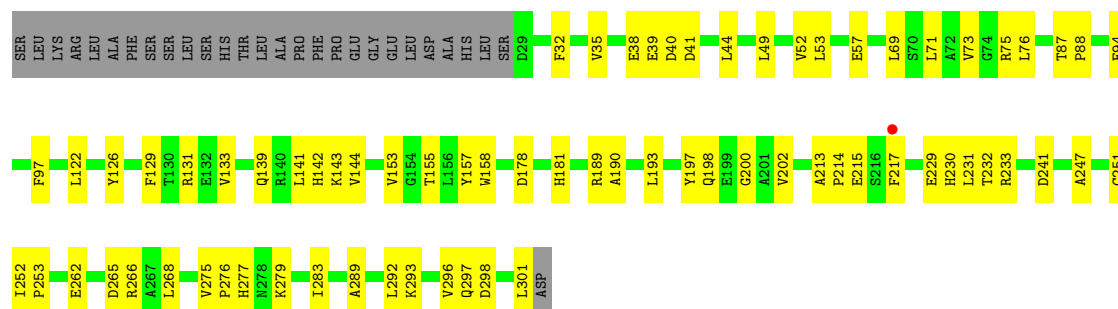
- Molecule 2 is a protein called TatT (Tp0956).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	273	Total	C	N	O	S	0	0	0
			2166	1386	374	398	8			
2	B	273	Total	C	N	O	S	0	0	0
			2166	1386	374	398	8			
2	C	273	Total	C	N	O	S	0	0	0
			2166	1386	374	398	8			

i

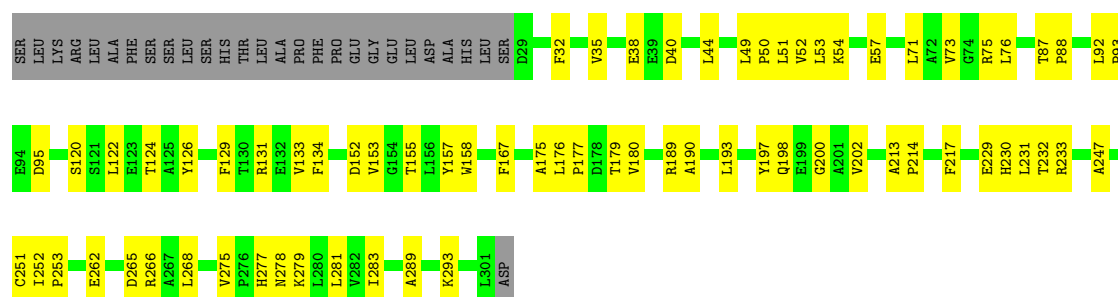
- Molecule 1: TatP(T) (Tp0957)





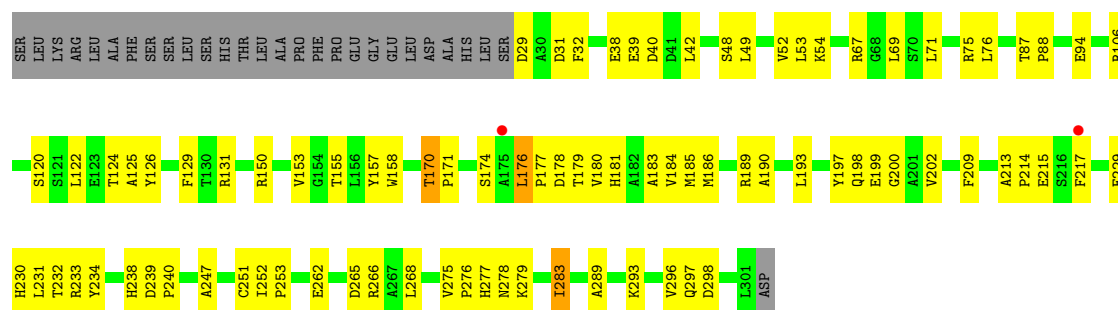
• Molecule 2: TatT (Tp0956)

Chain B: 67% 24% 9%



• Molecule 2: TatT (Tp0956)

Chain C: 62% 28% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	148.93Å 148.93Å 212.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.65 – 3.05 49.65 – 3.05	Depositor EDS
% Data completeness (in resolution range)	98.9 (49.65-3.05) 98.9 (49.65-3.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.203 , 0.263 (Not available) , 0.250	Depositor DCC
R_{free} test set	2321 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	97.6	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 100.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11376	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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 [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	D	0.54	0/2434	1.05	11/3292 (0.3%)
1	E	0.55	0/2546	0.98	7/3446 (0.2%)
2	A	0.59	0/2224	0.93	3/3029 (0.1%)
2	B	0.60	0/2224	0.93	2/3029 (0.1%)
2	C	0.60	0/2224	0.94	6/3029 (0.2%)
All	All	0.58	0/11652	0.97	29/15825 (0.2%)

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	D	0	1
1	E	0	2
All	All	0	3

There are no bond length outliers.

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	22	SER	N-CA-C	9.24	122.52	111.33
1	D	312	ASN	N-CA-C	8.97	123.54	112.23
2	A	275	VAL	CA-C-N	7.23	127.07	119.05
2	A	275	VAL	C-N-CA	7.23	127.07	119.05
2	C	176	LEU	CA-C-N	6.73	126.24	119.24

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	228	ILE	Peptide

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Mol	Chain	Res	Type	Group
1	E	228	ILE	Peptide
1	E	241	PRO	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	D	2385	0	2423	125	0
1	E	2493	0	2535	139	0
2	A	2166	0	2100	59	0
2	B	2166	0	2100	66	0
2	C	2166	0	2100	71	1
All	All	11376	0	11258	445	1

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 20.

The worst 5 of 445 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:E:240:ILE:HG21	1:E:248:MET:HE1	1.21	1.15
1:D:240:ILE:HG21	1:D:248:MET:HE1	1.17	1.15
1:D:104:ASN:HB3	1:D:107:ASP:HB2	1.23	1.14
1:E:65:SER:HA	1:E:66:SER:HB2	1.23	1.13
1:E:104:ASN:HB3	1:E:107:ASP:HB2	1.28	1.13

All (1) 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)
2:C:150:ARG:NH2	2:C:199:GLU:OE2[7_464]	2.12	0.08

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	D	297/324 (92%)	272 (92%)	24 (8%)	1 (0%)	36	63
1	E	284/324 (88%)	265 (93%)	16 (6%)	3 (1%)	11	34
2	A	271/301 (90%)	263 (97%)	8 (3%)	0	100	100
2	B	271/301 (90%)	258 (95%)	13 (5%)	0	100	100
2	C	271/301 (90%)	260 (96%)	11 (4%)	0	100	100
All	All	1394/1551 (90%)	1318 (94%)	72 (5%)	4 (0%)	36	63

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	23	ILE
1	E	242	SER
1	E	20	ALA
1	E	104	ASN

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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	D	303/324 (93%)	-0.20	2 (0%) 84 66	90, 129, 173, 200	0
1	E	317/324 (97%)	-0.11	2 (0%) 85 69	80, 128, 169, 210	0
2	A	273/301 (90%)	-0.55	1 (0%) 88 75	71, 103, 144, 177	0
2	B	273/301 (90%)	-0.44	0 100 100	66, 101, 146, 176	0
2	C	273/301 (90%)	-0.37	2 (0%) 84 66	73, 106, 147, 177	0
All	All	1439/1551 (92%)	-0.32	7 (0%) 87 72	66, 114, 163, 210	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	166	SER	3.3
1	E	187	LEU	3.2
2	C	217	PHE	2.8
2	A	217	PHE	2.6
1	D	167	VAL	2.2

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.