



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2024 – 09:40 AM EDT

PDB ID : 3WPE  
Title : Crystal structure of bovine TLR9 in complex with agonistic DNA1668\_12mer  
Authors : Ohto, U.; Tanji, H.; Shimizu, T.  
Deposited on : 2014-01-11  
Resolution : 2.38 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

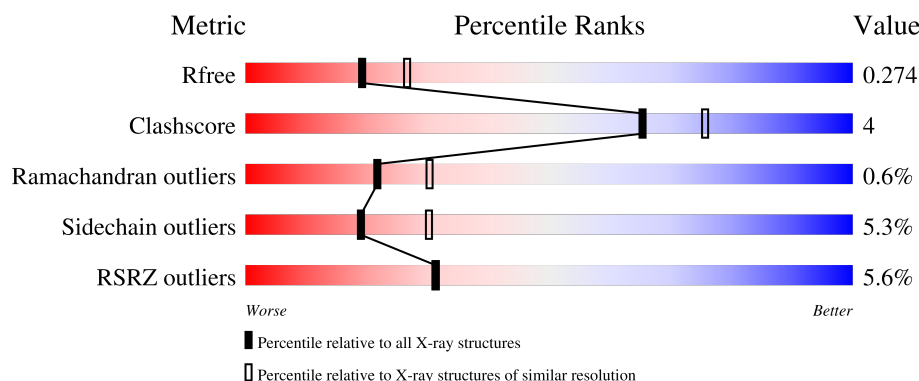
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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))
$R_{free}$	164625	6699 (2.40-2.36)
Clashscore	180529	7414 (2.40-2.36)
Ramachandran outliers	177936	7337 (2.40-2.36)
Sidechain outliers	177891	7338 (2.40-2.36)
RSRZ outliers	164620	6699 (2.40-2.36)

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  $\geq 3$ , 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	801	<div> <div>5%</div> <div> <div></div> <div>78%</div> <div>10%</div> <div>11%</div> </div> </div>
2	C	12	<div> <div>17%</div> <div>33%</div> <div>50%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5783 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 Toll-like receptor 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	714	Total	C	N	O	S	0	0	0
			5628	3596	996	1013	23			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	ARG	-	expression tag	UNP Q5I2M5
A	22	SER	-	expression tag	UNP Q5I2M5
A	23	PRO	-	expression tag	UNP Q5I2M5
A	24	TRP	-	expression tag	UNP Q5I2M5
A	816	GLU	-	expression tag	UNP Q5I2M5
A	817	PHE	-	expression tag	UNP Q5I2M5
A	818	LEU	-	expression tag	UNP Q5I2M5
A	819	VAL	-	expression tag	UNP Q5I2M5
A	820	PRO	-	expression tag	UNP Q5I2M5
A	821	ARG	-	expression tag	UNP Q5I2M5

- Molecule 2 is a DNA chain called DNA (5'-D(\*CP\*AP\*TP\*GP\*AP\*CP\*GP\*TP\*TP\*CP\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	P	0	0	0
			124	59	22	37	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		



- Molecule 1: Toll-like receptor 9



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	115.82Å 113.08Å 67.43Å 90.00° 100.61° 90.00°	Depositor
Resolution (Å)	43.05 – 2.38 43.05 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.05-2.38) 99.7 (43.05-2.38)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.217 , 0.274 0.218 , 0.274	Depositor DCC
$R_{free}$ test set	1717 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.8	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 43.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5783	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	A	0.33	0/5759	0.58	0/7827
2	C	0.29	0/138	0.69	0/211
All	All	0.33	0/5897	0.58	0/8038

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.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	469	CYS	Peptide

### 5.2 Too-close contacts [i](#)

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	5628	0	5661	40	0
2	C	124	0	69	4	0
3	A	31	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5783	0	5730	41	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 4.

All (41) 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:73:ARG:NH1	2:C:7:DG:O6	2.16	0.78
1:A:470:ASN:O	1:A:471:LEU:HD23	1.93	0.68
1:A:232:VAL:HG12	1:A:271:PHE:CD1	2.34	0.62
1:A:738:VAL:O	1:A:768:PHE:HB3	2.00	0.62
1:A:73:ARG:NH2	2:C:6:DC:OP2	2.36	0.59
1:A:275:HIS:HB2	1:A:276:PRO:HD2	1.85	0.58
1:A:642:LEU:HD13	1:A:644:ARG:HD2	1.85	0.58
1:A:390:PRO:O	1:A:415:SER:OG	2.18	0.58
1:A:379:THR:OG1	1:A:380:ASN:N	2.38	0.56
1:A:304:ARG:NH1	1:A:331:ASP:OD2	2.39	0.54
1:A:695:LEU:HD23	1:A:722:ALA:HB2	1.89	0.54
1:A:232:VAL:CG1	1:A:271:PHE:CD1	2.91	0.54
1:A:469:CYS:SG	1:A:469:CYS:O	2.66	0.53
1:A:44:CYS:HB2	1:A:69:LEU:HD23	1.91	0.52
1:A:295:LEU:H	1:A:317:ASN:HD22	1.59	0.51
1:A:73:ARG:HH21	2:C:5:DA:H3'	1.76	0.50
1:A:235:ALA:HB1	1:A:236:PRO:HD2	1.93	0.50
1:A:399:LEU:HD23	1:A:423:ASP:HB3	1.94	0.50
1:A:555:GLN:HB3	1:A:556:PRO:HD3	1.94	0.49
1:A:769:VAL:O	1:A:773:LEU:HD13	2.13	0.49
1:A:276:PRO:O	3:A:917:HOH:O	2.20	0.48
1:A:276:PRO:HB3	1:A:301:ASP:CB	2.43	0.48
1:A:293:SER:H	1:A:317:ASN:HD21	1.61	0.48
1:A:100:ALA:O	1:A:106:HIS:HA	2.14	0.47
1:A:359:PHE:O	1:A:362:LEU:HB2	2.15	0.46
1:A:194:LEU:O	1:A:218:PRO:HG3	2.14	0.46
1:A:49:LEU:H	1:A:72:ASN:HD22	1.66	0.44
2:C:4:DG:H2''	2:C:5:DA:C8	2.53	0.44
1:A:402:ILE:H	1:A:424:ASN:HD22	1.65	0.43
1:A:585:ILE:H	1:A:606:ASN:HD22	1.67	0.43
1:A:790:SER:HA	1:A:791:PRO:HA	1.89	0.43
1:A:775:ARG:HD2	1:A:775:ARG:HA	1.77	0.43
1:A:468:SER:OG	1:A:469:CYS:N	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:PHE:HA	1:A:372:GLY:O	2.19	0.42
1:A:761:HIS:ND1	1:A:790:SER:HB3	2.34	0.42
1:A:750:LEU:HD12	1:A:753:LEU:HB2	2.00	0.42
1:A:271:PHE:CE1	1:A:273:LYS:HD2	2.55	0.42
1:A:747:ALA:O	1:A:749:THR:N	2.53	0.42
1:A:728:LEU:HG	1:A:730:LEU:HD13	2.02	0.41
1:A:371:HIS:HB2	1:A:398:GLN:O	2.20	0.41
1:A:535:HIS:HE1	3:A:922:HOH:O	2.05	0.40

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	704/801 (88%)	653 (93%)	47 (7%)	4 (1%)	22	30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	720	VAL
1	A	361	SER
1	A	748	GLY
1	A	791	PRO

### 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	641/712 (90%)	607 (95%)	34 (5%)	19	30

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	ARG
1	A	119	LEU
1	A	183	CYS
1	A	245	LEU
1	A	265	ARG
1	A	269	LYS
1	A	308	ARG
1	A	377	SER
1	A	379	THR
1	A	385	SER
1	A	388	ARG
1	A	469	CYS
1	A	470	ASN
1	A	471	LEU
1	A	492	ARG
1	A	544	GLN
1	A	601	LEU
1	A	619	LEU
1	A	623	LYS
1	A	640	THR
1	A	649	LEU
1	A	653	LEU
1	A	671	LEU
1	A	673	VAL
1	A	687	LEU
1	A	695	LEU
1	A	700	ARG
1	A	721	ARG
1	A	735	LEU
1	A	743	PHE
1	A	749	THR
1	A	750	LEU
1	A	771	PHE
1	A	790	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	132	ASN
1	A	317	ASN
1	A	330	ASN
1	A	334	GLN
1	A	396	HIS
1	A	400	ASN
1	A	424	ASN
1	A	544	GLN
1	A	583	ASN
1	A	606	ASN
1	A	627	ASN
1	A	702	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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	714/801 (89%)	0.39	40 (5%) 31 31	23, 47, 78, 111	0
2	C	6/12 (50%)	0.01	0 100 100	54, 56, 60, 65	0
All	All	720/813 (88%)	0.38	40 (5%) 31 31	23, 47, 78, 111	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	747	ALA	4.6
1	A	769	VAL	4.2
1	A	29	PRO	4.1
1	A	763	ALA	3.8
1	A	773	LEU	3.7
1	A	269	LYS	3.6
1	A	469	CYS	3.5
1	A	744	GLY	3.4
1	A	361	SER	3.1
1	A	36	LEU	3.0
1	A	767	ALA	3.0
1	A	792	GLY	3.0
1	A	779	VAL	3.0
1	A	775	ARG	3.0
1	A	772	LEU	2.9
1	A	533	LEU	2.8
1	A	468	SER	2.8
1	A	780	PRO	2.8
1	A	534	TYR	2.8
1	A	768	PHE	2.7
1	A	185	ARG	2.7
1	A	672	THR	2.6
1	A	270	ASN	2.6
1	A	184	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	750	LEU	2.4
1	A	39	HIS	2.3
1	A	304	ARG	2.3
1	A	791	PRO	2.2
1	A	275	HIS	2.2
1	A	352	HIS	2.2
1	A	786	VAL	2.2
1	A	64	VAL	2.2
1	A	370	MET	2.2
1	A	30	ALA	2.1
1	A	749	THR	2.1
1	A	778	ALA	2.1
1	A	296	TYR	2.1
1	A	38	PRO	2.1
1	A	276	PRO	2.0
1	A	183	CYS	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 monosaccharides 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.