



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:16 PM GMT

PDB ID : 4Y04  
Title : Crystal structure of dipeptidyl peptidase 11 (DPP11) from Porphyromonas gingivalis (Space)  
Authors : Sakamoto, Y.; Suzuki, Y.; Iizuka, I.; Tateoka, C.; Roppongi, S.; Fujimoto, M.; Inaka, K.; Tanaka, H.; Yamada, M.; Ohta, K.; Nonaka, T.; Ogasawara, W.; Tanaka, N.  
Deposited on : 2015-02-05  
Resolution : 1.66 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

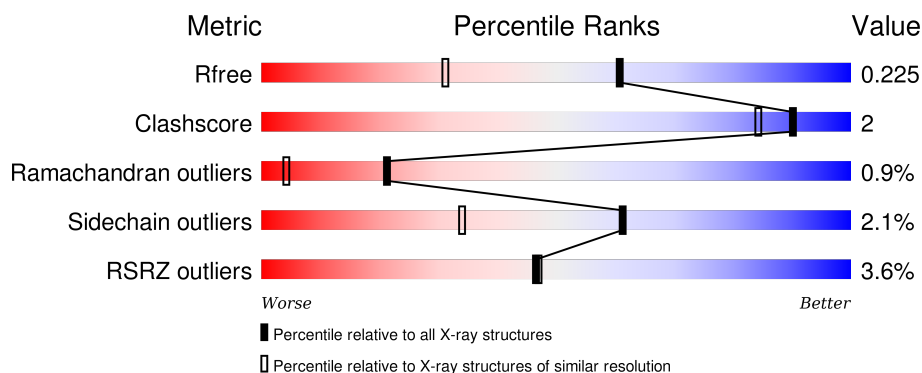
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 1.66 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))
$R_{free}$	91344	1226 (1.66-1.66)
Clashscore	102246	1323 (1.66-1.66)
Ramachandran outliers	100387	1295 (1.66-1.66)
Sidechain outliers	100360	1295 (1.66-1.66)
RSRZ outliers	91569	1227 (1.66-1.66)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	720	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6237 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 Peptidase S46.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	687	Total	C	N	O	S	0	1	0
			5533	3510	959	1037	27			

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	K	0	0
			2	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



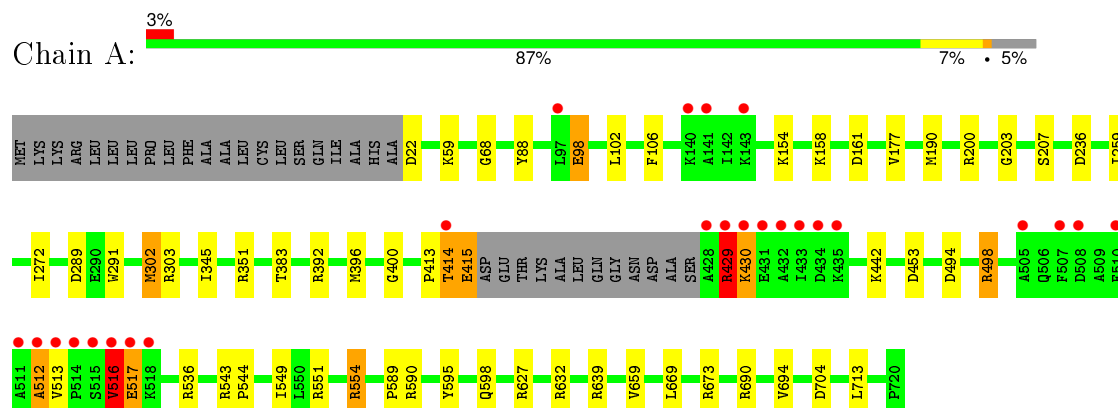
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	696	Total 696	O 696	0	0



- Molecule 1: Peptidase S46



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	99.14Å 103.28Å 175.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.66 49.57 – 1.66	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-1.66) 99.9 (49.57-1.66)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.49 (at 1.66Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.186 , 0.225 0.193 , 0.225	Depositor DCC
$R_{free}$ test set	5266 reflections (5.23%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.1	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 37.1	EDS
Estimated twinning fraction	0.058 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.24$	Xtriage
Outliers	0 of 105865 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6237	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, K

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	1.01	0/5657	1.06	29/7640 (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	3

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392	ARG	NE-CZ-NH2	-14.05	113.27	120.30
1	A	392	ARG	NE-CZ-NH1	11.35	125.97	120.30
1	A	590	ARG	NE-CZ-NH1	9.95	125.27	120.30
1	A	498	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	551	ARG	NE-CZ-NH2	-6.88	116.86	120.30
1	A	690	ARG	NE-CZ-NH1	6.74	123.67	120.30
1	A	396	MET	CG-SD-CE	-6.68	89.51	100.20
1	A	632	ARG	NE-CZ-NH1	6.55	123.58	120.30
1	A	551	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	A	351	ARG	NE-CZ-NH2	-6.38	117.11	120.30
1	A	590	ARG	NE-CZ-NH2	-6.21	117.19	120.30
1	A	351	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	A	161	ASP	CB-CG-OD1	6.17	123.85	118.30
1	A	453	ASP	CB-CG-OD2	-6.15	112.77	118.30
1	A	303	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	A	414	THR	N-CA-CB	5.67	121.07	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	516	VAL	N-CA-C	5.51	125.88	111.00
1	A	200	ARG	NE-CZ-NH1	5.33	122.96	120.30
1	A	512	ALA	N-CA-C	5.30	125.32	111.00
1	A	236	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	303	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	536	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	498	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	289	ASP	CB-CG-OD1	5.18	122.96	118.30
1	A	302	MET	CG-SD-CE	5.17	108.47	100.20
1	A	554	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	A	516	VAL	C-N-CA	5.10	134.46	121.70
1	A	536	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	516	VAL	CA-C-N	5.07	128.36	117.20

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	413	PRO	Peptide
1	A	429	ARG	Peptide
1	A	516	VAL	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	5533	0	5424	25	0
2	A	2	0	0	0	0
3	A	6	0	8	0	0
4	A	696	0	0	3	1
All	All	6237	0	5432	25	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 2.

All (25) 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:673[A]:ARG:NH1	4:A:1588:HOH:O	1.88	0.88
1:A:302:MET:HE2	1:A:400:GLY:HA2	1.58	0.84
1:A:516:VAL:N	1:A:517:GLU:HB3	2.06	0.71
1:A:669:LEU:HB3	1:A:694:VAL:HG12	1.77	0.66
1:A:429:ARG:CB	1:A:430:LYS:HB3	2.28	0.63
1:A:627:ARG:HD3	4:A:912:HOH:O	2.01	0.59
1:A:429:ARG:HB2	1:A:430:LYS:HB3	1.87	0.57
1:A:669:LEU:HB3	1:A:694:VAL:CG1	2.35	0.57
1:A:414:THR:HA	1:A:415:GLU:HB3	1.91	0.52
1:A:554:ARG:NH1	4:A:1280:HOH:O	2.42	0.51
1:A:494:ASP:OD1	1:A:498:ARG:HD3	2.12	0.50
1:A:88:TYR:CE1	1:A:102:LEU:HD11	2.48	0.48
1:A:302:MET:CE	1:A:400:GLY:HA2	2.39	0.48
1:A:429:ARG:CA	1:A:430:LYS:HB3	2.45	0.46
1:A:383:THR:HG21	1:A:549:ILE:HG13	1.98	0.46
1:A:177:VAL:HG11	1:A:190:MET:HE2	1.98	0.45
1:A:589:PRO:HG3	1:A:595:TYR:CE2	2.53	0.44
1:A:494:ASP:OD2	1:A:498:ARG:NH1	2.43	0.44
1:A:704:ASP:HB2	1:A:713:LEU:HD11	1.99	0.44
1:A:177:VAL:CG1	1:A:190:MET:HE2	2.48	0.44
1:A:543:ARG:HB3	1:A:544:PRO:HD3	1.99	0.43
1:A:272:ILE:HG12	1:A:659:VAL:HG22	2.01	0.42
1:A:177:VAL:HB	1:A:190:MET:CE	2.50	0.42
1:A:106:PHE:O	1:A:203:GLY:HA2	2.21	0.41
1:A:291:TRP:CE3	1:A:345:ILE:HD11	2.56	0.40

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)
4:A:981:HOH:O	4:A:981:HOH:O[3_555]	1.81	0.39

## 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	684/720 (95%)	664 (97%)	14 (2%)	6 (1%)	21 4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	GLU
1	A	512	ALA
1	A	517	GLU
1	A	68	GLY
1	A	430	LYS
1	A	513	VAL

### 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	582/607 (96%)	570 (98%)	12 (2%)	61 34

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	59	LYS
1	A	98	GLU
1	A	154	LYS
1	A	158	LYS
1	A	207	SER
1	A	259	ILE
1	A	415	GLU
1	A	429	ARG
1	A	442	LYS
1	A	598	GLN
1	A	639	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	307	GLN
1	A	353	ASN
1	A	378	HIS
1	A	506	GLN
1	A	569	GLN
1	A	654	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z  > 2$
3	GOL	A	803	-	5,5,5	0.52	0	5,5,5	0.47	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	803	-	-	0/4/4/4	0/0/0/0

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, 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	687/720 (95%)	-0.20	25 (3%)	46 47	19, 29, 53, 95	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	97	LEU	8.2
1	A	516	VAL	7.4
1	A	513	VAL	5.2
1	A	432	ALA	4.8
1	A	511	ALA	4.6
1	A	433	ILE	4.1
1	A	510	PHE	4.1
1	A	430	LYS	3.9
1	A	512	ALA	3.9
1	A	515	SER	3.6
1	A	429	ARG	3.5
1	A	414	THR	3.2
1	A	518	LYS	3.2
1	A	514	PRO	3.0
1	A	517	GLU	2.9
1	A	143	LYS	2.9
1	A	428	ALA	2.7
1	A	140	LYS	2.7
1	A	431	GLU	2.5
1	A	434	ASP	2.5
1	A	141	ALA	2.3
1	A	435	LYS	2.3
1	A	507	PHE	2.2
1	A	508	ASP	2.1
1	A	505	ALA	2.1

## 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	A	803	6/6	0.95	0.08	-1.14	28,33,34,37	0
2	K	A	802	1/1	0.99	0.06	-1.23	28,28,28,28	0
2	K	A	801	1/1	0.98	0.05	-	31,31,31,31	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.