



# wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 11:37 PM GMT

PDB ID : 3PJI  
Title : Crystal structure of AgamOBP22a at 1.7 angstrom in the open status for ligand binding.  
Authors : Zhang, S.; Ren, H.  
Deposited on : 2010-11-10  
Resolution : 1.70 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

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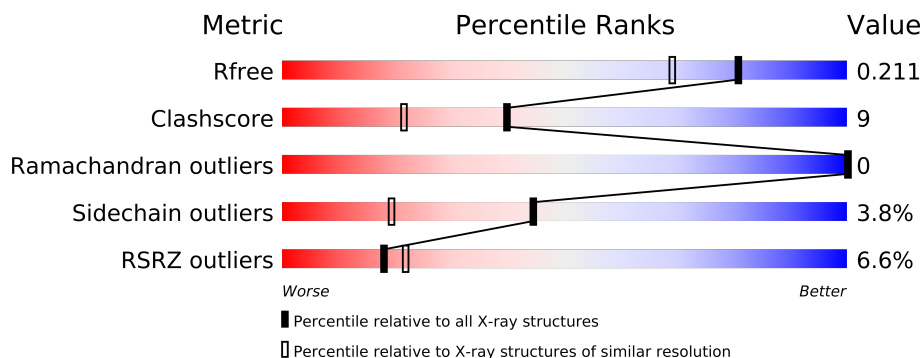
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.70 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))
$R_{free}$	66092	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	141	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NH4	A	202	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1115 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Odorant binding protein (AGAP010409-PA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	6	0
			994	618	164	205	7			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	LEU	-	EXPRESSION TAG	UNP Q7PGA3
A	146	VAL	-	EXPRESSION TAG	UNP Q7PGA3
A	147	PRO	-	EXPRESSION TAG	UNP Q7PGA3
A	148	ARG	-	EXPRESSION TAG	UNP Q7PGA3
A	149	GLY	-	EXPRESSION TAG	UNP Q7PGA3
A	150	SER	-	EXPRESSION TAG	UNP Q7PGA3
A	151	SER	-	EXPRESSION TAG	UNP Q7PGA3
A	152	GLY	-	EXPRESSION TAG	UNP Q7PGA3
A	153	GLU	-	EXPRESSION TAG	UNP Q7PGA3
A	154	LEU	-	EXPRESSION TAG	UNP Q7PGA3
A	155	TRP	-	EXPRESSION TAG	UNP Q7PGA3
A	156	SER	-	EXPRESSION TAG	UNP Q7PGA3
A	157	HIS	-	EXPRESSION TAG	UNP Q7PGA3
A	158	PRO	-	EXPRESSION TAG	UNP Q7PGA3
A	159	GLN	-	EXPRESSION TAG	UNP Q7PGA3
A	160	PHE	-	EXPRESSION TAG	UNP Q7PGA3
A	161	GLU	-	EXPRESSION TAG	UNP Q7PGA3
A	162	LYS	-	EXPRESSION TAG	UNP Q7PGA3

- Molecule 2 is AMMONIUM ION (three-letter code: NH<sub>4</sub>) (formula: H<sub>4</sub>N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0
2	A	1	Total N 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	118	Total O 118 118	0	0

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- Molecule 1: Odorant binding protein (AGAP010409-PA)

Amino Acid	Relative Abundance
PHE	125
ASP	126
ASN	127
ALA	128
THR	129
LEU	130
VAL	131
PRO	132
ARG	133
GLY	134
SER	135
CYS	136
LYS	137
GLU	138

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.26Å 37.70Å 43.64Å 90.00° 99.46° 90.00°	Depositor
Resolution (Å)	25.42 – 1.70 25.42 – 1.70	Depositor EDS
% Data completeness (in resolution range)	88.9 (25.42-1.70) 91.9 (25.42-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.59 (at 1.71Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.179 , 0.208 0.181 , 0.211	Depositor DCC
$R_{free}$ test set	1163 reflections (9.86%)	DCC
Wilson B-factor (Å <sup>2</sup> )	16.4	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 39.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 11791 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	1115	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.47% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NH4

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	0/1035	0.50	0/1402

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	994	0	895	18	0
2	A	3	0	0	1	0
3	A	118	0	0	4	0
All	All	1115	0	895	18	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

The worst 5 of 18 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:96[A]:GLU:HG3	2:A:200:NH4:N	1.94	0.81

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:112:GLU:HB2	3:A:292:HOH:O	1.95	0.65
1:A:28:VAL:HG12	1:A:66:ILE:CD1	2.27	0.64
1:A:153:GLU:N	3:A:275:HOH:O	2.32	0.61
1:A:28:VAL:HG12	1:A:66:ILE:HD13	1.86	0.57

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	121/141 (86%)	120 (99%)	1 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	110/122 (90%)	106 (96%)	4 (4%)	47	22

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	107	GLU
1	A	112	GLU
1	A	115	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such



sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	90	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 are modelled with single atom - leaving 0 for Mogul analysis.

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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	119/141 (84%)	0.10	8 (6%) 17 21	9, 17, 36, 49	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	158	PRO	5.1
1	A	125[A]	PHE	3.6
1	A	137	ASN	3.6
1	A	25	ASN	2.9
1	A	135	TYR	2.3

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NH4	A	202	1/1	0.12	3.11	12,12,12,12	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NH4	A	200	1/1	0.10	0.27	13,13,13,13	0
2	NH4	A	201	1/1	0.09	-0.63	11,11,11,11	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.