



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:04 PM GMT

PDB ID : 4J6T  
Title : Crystal Structure of Tyrosinase from Bacillus megaterium F197A mutant  
Authors : Kanteev, M.; Goldfeder, M.; Adir, N.; Fishman, A.  
Deposited on : 2013-02-12  
Resolution : 2.43 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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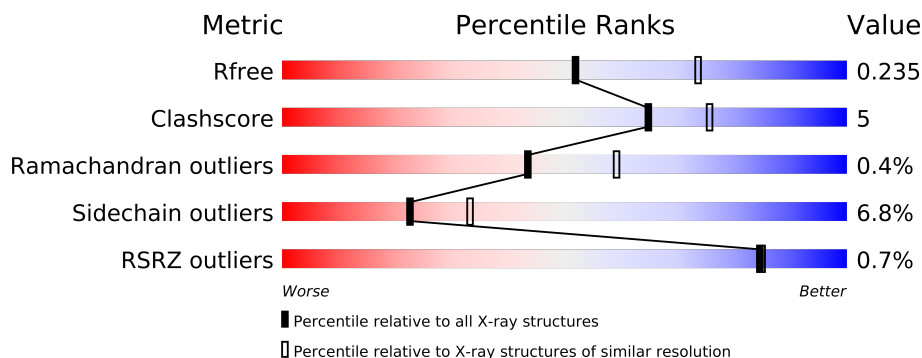
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.43 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	2989 (2.48-2.40)
Clashscore	79885	3698 (2.48-2.40)
Ramachandran outliers	78287	3639 (2.48-2.40)
Sidechain outliers	78261	3640 (2.48-2.40)
RSRZ outliers	66119	2993 (2.48-2.40)

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	303	
1	B	303	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4899 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 Tyrosinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2348	1492	424	424	8			
1	B	286	Total	C	N	O	S	0	0	0
			2340	1486	423	423	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	SER	ENGINEERED MUTATION	UNP B2ZB02
A	197	ALA	PHE	ENGINEERED MUTATION	UNP B2ZB02
A	298	HIS	-	EXPRESSION TAG	UNP B2ZB02
A	299	HIS	-	EXPRESSION TAG	UNP B2ZB02
A	300	HIS	-	EXPRESSION TAG	UNP B2ZB02
A	301	HIS	-	EXPRESSION TAG	UNP B2ZB02
A	302	HIS	-	EXPRESSION TAG	UNP B2ZB02
A	303	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	2	GLY	SER	ENGINEERED MUTATION	UNP B2ZB02
B	197	ALA	PHE	ENGINEERED MUTATION	UNP B2ZB02
B	298	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	299	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	300	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	301	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	302	HIS	-	EXPRESSION TAG	UNP B2ZB02
B	303	HIS	-	EXPRESSION TAG	UNP B2ZB02

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Cu	0	0
			2	2		
2	A	2	Total	Cu	0	0
			2	2		

- Molecule 3 is water.

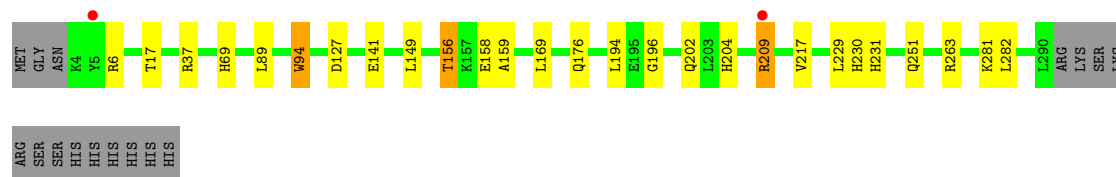
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	121	Total 121	O 121	0	0
3	B	86	Total 86	O 86	0	0

### 3 Residue-property plots i

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

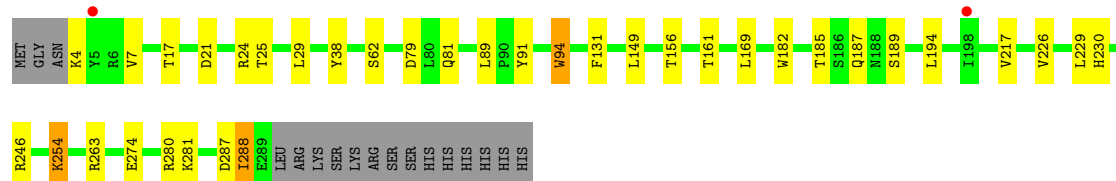
#### • Molecule 1: Tyrosinase

Chain A: 



#### • Molecule 1: Tyrosinase

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.01Å 78.81Å 84.34Å 90.00° 106.62° 90.00°	Depositor
Resolution (Å)	27.48 – 2.43 27.48 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.1 (27.48-2.43) 98.9 (27.48-2.43)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.76 (at 2.42Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, $R_{free}$	0.196 , 0.237 0.193 , 0.235	Depositor DCC
$R_{free}$ test set	1316 reflections (5.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.0	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 20.5	EDS
Estimated twinning fraction	0.034 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 25963 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4899	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.34% 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: CU

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	0.50	0/2425	0.59	0/3305
1	B	0.45	0/2417	0.58	0/3294
All	All	0.48	0/4842	0.58	0/6599

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	2348	0	0	11	0
1	B	2340	0	0	14	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	121	0	0	3	0
3	B	86	0	0	3	0
All	All	4899	0	0	25	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 5.

All (25) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:159:ALA:CB	1:A:209:ARG:CG	2.46	0.94
1:A:156:THR:CG2	1:A:209:ARG:NE	2.44	0.81
1:B:287:ASP:O	1:B:288:ILE:CB	2.35	0.75
1:B:287:ASP:O	1:B:288:ILE:CG2	2.39	0.70
1:A:196:GLY:CA	1:A:202:GLN:O	2.49	0.61
1:B:62:SER:OG	1:B:182:TRP:O	2.24	0.56
1:B:38:TYR:OH	1:B:79:ASP:OD1	2.25	0.55
1:B:21:ASP:O	1:B:25:THR:OG1	2.26	0.54
1:B:91:TYR:CE2	1:B:287:ASP:OD1	2.61	0.54
1:B:263:ARG:NH1	1:B:274:GLU:OE1	2.42	0.53
1:A:204:HIS:CD2	1:A:230:HIS:CE1	2.97	0.52
1:A:176:GLN:OE1	1:A:251:GLN:NE2	2.43	0.51
1:B:189:SER:N	3:B:564:HOH:O	2.44	0.51
1:B:91:TYR:OH	1:B:287:ASP:OD1	2.30	0.50
1:A:202:GLN:NE2	3:A:607:HOH:O	2.44	0.50
1:A:69:HIS:CD2	1:A:231:HIS:CE1	3.01	0.49
1:B:131:PHE:N	3:B:505:HOH:O	2.46	0.49
1:A:94:TRP:NE1	1:A:230:HIS:CD2	2.80	0.48
1:B:287:ASP:OD1	1:B:287:ASP:O	2.31	0.48
1:B:94:TRP:NE1	1:B:230:HIS:CD2	2.82	0.47
1:A:263:ARG:NE	3:A:561:HOH:O	2.46	0.47
1:A:37:ARG:NH1	3:A:582:HOH:O	2.48	0.46
1:B:187:GLN:N	3:B:559:HOH:O	2.48	0.46
1:A:6:ARG:NH2	1:A:282:LEU:O	2.50	0.45
1:B:185:THR:O	1:B:187:GLN:NE2	2.53	0.42

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	285/303 (94%)	274 (96%)	11 (4%)	0	100	100
1	B	284/303 (94%)	272 (96%)	10 (4%)	2 (1%)	30	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	569/606 (94%)	546 (96%)	21 (4%)	2 (0%)	43	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	288	ILE
1	B	254	LYS

### 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	249/264 (94%)	235 (94%)	14 (6%)	30	44
1	B	248/264 (94%)	228 (92%)	20 (8%)	17	24
All	All	497/528 (94%)	463 (93%)	34 (7%)	22	33

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

Mol	Chain	Res	Type
1	A	17	THR
1	A	89	LEU
1	A	94	TRP
1	A	127	ASP
1	A	141	GLU
1	A	149	LEU
1	A	156	THR
1	A	158	GLU
1	A	169	LEU
1	A	194	LEU
1	A	209	ARG
1	A	217	VAL
1	A	229	LEU
1	A	281	LYS
1	B	4	LYS
1	B	7	VAL
1	B	17	THR

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Mol	Chain	Res	Type
1	B	24	ARG
1	B	29	LEU
1	B	81	GLN
1	B	89	LEU
1	B	94	TRP
1	B	149	LEU
1	B	156	THR
1	B	161	THR
1	B	169	LEU
1	B	194	LEU
1	B	217	VAL
1	B	226	VAL
1	B	229	LEU
1	B	246	ARG
1	B	254	LYS
1	B	280	ARG
1	B	281	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 4 ligands modelled in this entry, 4 are monoatomic - 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	287/303 (94%)	-0.24	2 (0%) 84 85	17, 27, 48, 87	1 (0%)
1	B	286/303 (94%)	-0.19	2 (0%) 84 85	18, 34, 56, 95	0
All	All	573/606 (94%)	-0.22	4 (0%) 84 85	17, 31, 54, 95	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	ILE	3.4
1	B	5	TYR	2.6
1	A	209	ARG	2.5
1	A	5	TYR	2.2

### 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( $\text{\AA}^2$ )	Q<0.9
2	CU	B	402	1/1	0.14	-0.86	68,68,68,68	1
2	CU	B	401	1/1	0.05	-2.60	40,40,40,40	1
2	CU	A	401	1/1	0.05	-3.87	35,35,35,35	1
2	CU	A	402	1/1	0.06	-4.17	58,58,58,58	1

## 6.5 Other polymers ⓘ

There are no such residues in this entry.