



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:00 PM GMT

PDB ID : 3KVC  
Title : Crystal structure of bovine RPE65 at 1.9 angstrom resolution  
Authors : Kiser, P.D.; Lodowski, D.T.; Golczak, M.; Palczewski, K.  
Deposited on : 2009-11-29  
Resolution : 1.90 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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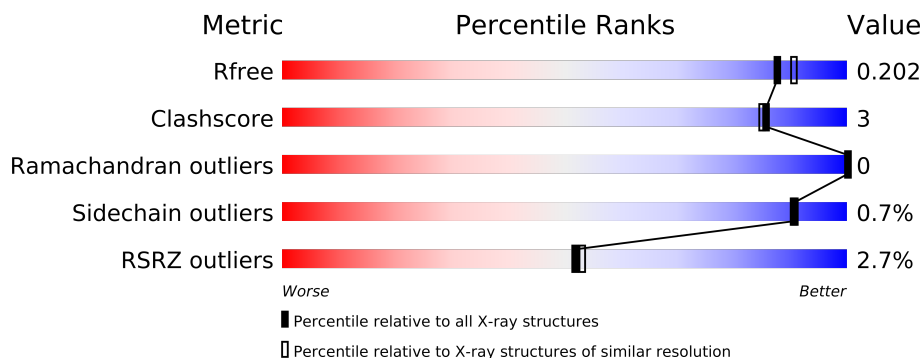
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.90 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	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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	533	
1	B	533	

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	FE2	B	534	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8556 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 Retinoid isomerohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	0	4	0
			4024	2597	663	748	16			
1	B	497	Total	C	N	O	S	0	5	0
			4027	2600	664	747	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	341	LEU	SER	VARIANT	UNP Q28175
B	341	LEU	SER	VARIANT	UNP Q28175

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	237	Total	O	0	0
			237	237		
3	B	266	Total	O	0	0
			266	266		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	176.36Å 176.36Å 86.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.90 50.91 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (20.00-1.90) 99.3 (50.91-1.90)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.75 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.146 , 0.171 0.177 , 0.202	Depositor DCC
$R_{free}$ test set	6166 reflections (5.13%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.9	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 36.4	EDS
Estimated twinning fraction	0.665 for H, K, L 0.335 for H+K, -K, -L 0.022 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.665 for H, K, L 0.335 for H+K, -K, -L	Depositor
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 120467 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8556	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: FE2

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	0.74	0/4148	0.72	2/5645 (0.0%)
1	B	0.77	0/4155	0.73	0/5655
All	All	0.76	0/8303	0.72	2/11300 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ASP	CB-CG-OD1	5.52	123.27	118.30
1	A	26	LEU	CA-CB-CG	5.13	127.10	115.30

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	4024	0	3928	21	0
1	B	4027	0	3932	23	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	237	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	266	0	0	1	0
All	All	8556	0	7860	43	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 3.

All (43) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:291:ILE:HD13	1:B:324:LEU:HD12	1.53	0.91
1:A:195[B]:CYS:SG	1:A:196:PHE:N	2.48	0.86
1:A:247:PRO:O	1:A:294:LYS:HE3	1.91	0.70
1:A:321:HIS:O	3:A:636:HOH:O	2.14	0.66
1:B:291:ILE:CD1	1:B:324:LEU:HD12	2.29	0.60
1:B:322:GLU:OE2	1:B:322:GLU:HA	2.01	0.60
1:B:141:GLU:HG3	1:B:212:LEU:HD22	1.84	0.59
1:B:394:ILE:HD12	3:B:737:HOH:O	2.01	0.59
1:A:207:VAL:HG22	1:A:225:ILE:HD13	1.86	0.58
1:B:291:ILE:HD13	1:B:324:LEU:CD1	2.32	0.57
1:B:215:ASP:OD2	1:B:215:ASP:N	2.38	0.56
1:B:491:VAL:HG12	1:B:493:PRO:HD3	1.88	0.56
1:A:26:LEU:C	1:A:26:LEU:HD23	2.28	0.54
1:B:412:PRO:O	1:B:414[A]:GLN:HG2	2.10	0.51
1:B:291:ILE:HD11	1:B:370:LEU:HD23	1.93	0.51
1:A:310:ASN:HB2	1:A:331:TRP:HB2	1.92	0.51
1:B:255:THR:HB	1:B:256:PRO:HD2	1.92	0.50
1:A:245:LEU:HD11	1:A:294:LYS:HD3	1.93	0.50
1:B:212:LEU:HD11	1:B:216:LYS:HA	1.93	0.50
1:A:214:ALA:O	1:A:216:LYS:HG2	2.12	0.50
1:B:310:ASN:HB2	1:B:331:TRP:HB2	1.94	0.49
1:A:307:SER:HB2	1:A:308:PRO:HD2	1.95	0.48
1:A:255:THR:HB	1:A:256:PRO:HD2	1.93	0.48
1:B:415:ALA:O	1:B:438:GLY:HA2	2.14	0.48
1:A:394:ILE:HD11	1:B:333:GLY:HA2	1.96	0.47
1:A:202:ILE:HG21	1:A:237:PRO:HB3	1.97	0.46
1:B:529:LEU:C	1:B:529:LEU:HD12	2.36	0.46
1:B:307:SER:HB2	1:B:308:PRO:HD2	1.97	0.45
1:A:183:ILE:HD12	1:A:189:VAL:HG22	1.99	0.45
1:A:34:ILE:HG12	1:A:504:ILE:HD12	1.98	0.45
1:A:47:PRO:HD2	1:A:524:VAL:HG13	1.98	0.45
1:A:276:MET:HG2	1:A:342:TYR:OH	2.17	0.44
1:B:460:TRP:CZ2	1:B:462:GLU:HG3	2.53	0.44
1:A:412:PRO:O	1:A:414[A]:GLN:HG2	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:138:PRO:HD2	1:B:183:ILE:CD1	2.48	0.44
1:A:102:GLU:O	1:A:103:PHE:C	2.55	0.44
1:A:529:LEU:C	1:A:529:LEU:HD12	2.39	0.43
1:B:246:THR:HB	1:B:247:PRO:HD2	2.01	0.42
1:A:291:ILE:HD13	1:A:291:ILE:HG21	1.81	0.41
1:B:426:TYR:CE1	1:B:481:GLU:HG3	2.56	0.41
1:A:207:VAL:HG22	1:A:225:ILE:CD1	2.50	0.41
1:B:89:TYR:O	1:B:93:MET:HG2	2.21	0.41
1:B:505:LEU:HD23	1:B:512:GLU:HA	2.02	0.41

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	494/533 (93%)	479 (97%)	15 (3%)	0	100	100
1	B	494/533 (93%)	482 (98%)	12 (2%)	0	100	100
All	All	988/1066 (93%)	961 (97%)	27 (3%)	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	445/475 (94%)	443 (100%)	2 (0%)	95	95
1	B	446/475 (94%)	442 (99%)	4 (1%)	87	86
All	All	891/950 (94%)	885 (99%)	6 (1%)	91	90



All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	215	ASP
1	B	201	SER
1	B	202	ILE
1	B	215	ASP
1	B	310	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	373	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 2 ligands modelled in this entry, 2 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	498/533 (93%)	-0.13	12 (2%) 56 57	18, 28, 43, 53	0
1	B	497/533 (93%)	-0.15	14 (2%) 50 51	17, 27, 41, 54	0
All	All	995/1066 (93%)	-0.14	26 (2%) 52 55	17, 28, 42, 54	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	496	GLY	4.9
1	B	496	GLY	4.9
1	A	272	GLY	4.9
1	B	196	PHE	4.8
1	A	3	SER	4.7
1	A	495	ALA	4.7
1	B	3	SER	4.4
1	B	197	GLY	4.1
1	B	495	ALA	3.8
1	A	197	GLY	3.8
1	B	214	ALA	3.7
1	B	235	PHE	3.3
1	A	214	ALA	2.8
1	A	273	ALA	2.8
1	B	273	ALA	2.8
1	A	4	GLN	2.7
1	B	497	GLN	2.6
1	A	196	PHE	2.6
1	B	201	SER	2.5
1	A	497	GLN	2.3
1	A	33	ARG	2.3
1	B	533	SER	2.3
1	B	234	ARG	2.2
1	A	201	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	202	ILE	2.1
1	B	477	ASP	2.1

## 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	FE2	B	534	1/1	0.14	3.29	21,21,21,21	0
2	FE2	A	534	1/1	0.10	-0.53	22,22,22,22	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.