



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:54 AM GMT

PDB ID : 3DPT  
Title : COR domain of Rab family protein (Roco)  
Authors : Gotthardt, K.; Weyand, M.; Kortholt, A.; Van Haastert, P.J.M.; Wittinghofer, A.  
Deposited on : 2008-07-09  
Resolution : 2.90 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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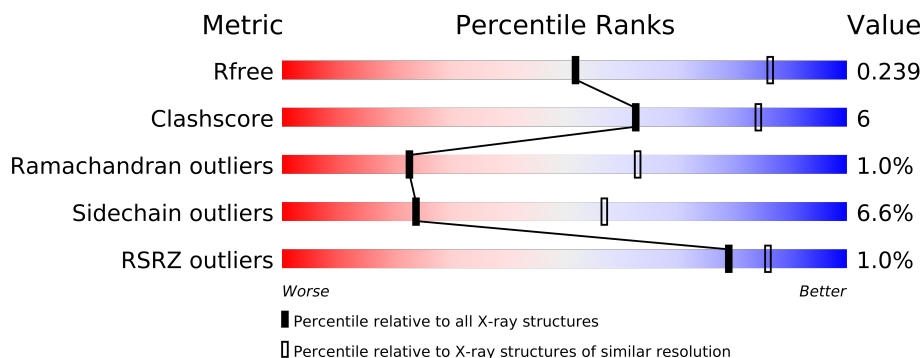
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.90 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	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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	332	
1	B	332	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4643 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 Rab family protein.

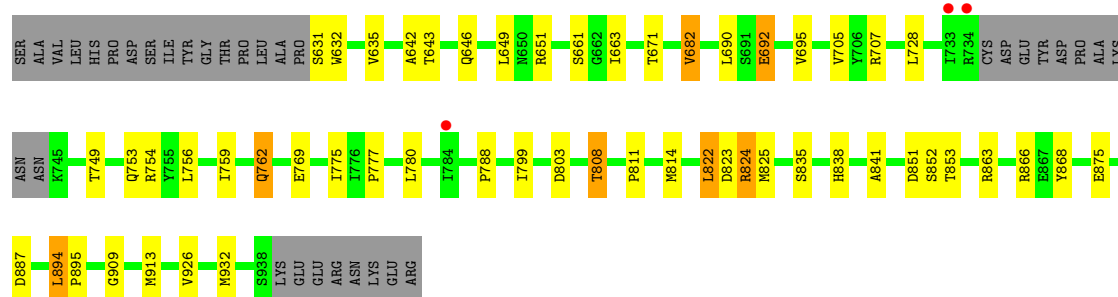
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	0
			2329	1502	373	443	11			
1	B	299	Total	C	N	O	S	0	0	0
			2314	1489	367	447	11			

### 3 Residue-property plots

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.

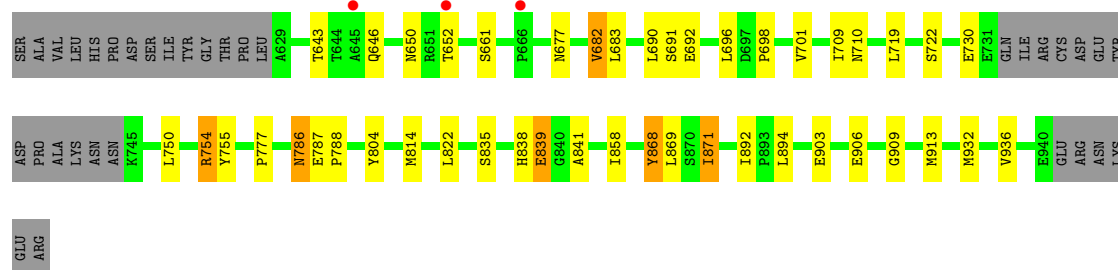
- Molecule 1: Rab family protein

Chain A: 



- Molecule 1: Rab family protein

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.11Å 130.11Å 129.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.83 – 2.90 45.84 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.83-2.90) 99.9 (45.84-2.90)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.79 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.210 , 0.243 0.211 , 0.239	Depositor DCC
$R_{free}$ test set	1422 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	66.7	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 58.5	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 28439 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4643	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% 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

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	0.44	0/2375	0.57	0/3231
1	B	0.43	0/2362	0.58	0/3218
All	All	0.43	0/4737	0.57	0/6449

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	2329	0	2261	30	0
1	B	2314	0	2195	23	0
All	All	4643	0	4456	51	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 6.

All (51) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:786:ASN:HD22	1:B:786:ASN:H	1.17	0.90
1:A:631:SER:O	1:A:635:VAL:HG23	1.81	0.81

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:635:VAL:HG22	1:A:661:SER:HB2	1.69	0.75
1:A:635:VAL:HG21	1:A:663:ILE:HD12	1.69	0.73
1:B:696:LEU:HD11	1:B:755:TYR:HE2	1.61	0.65
1:A:682:VAL:HG13	1:A:695:VAL:HG13	1.78	0.65
1:A:707:ARG:NH1	1:A:728:LEU:O	2.32	0.63
1:A:780:LEU:HD11	1:A:814:MET:HE1	1.82	0.62
1:A:823:ASP:H	1:A:825:MET:HG3	1.64	0.62
1:A:754:ARG:CZ	1:B:871:ILE:HD12	2.29	0.62
1:A:635:VAL:HG22	1:A:661:SER:CB	2.30	0.61
1:A:835:SER:HB2	1:A:838:HIS:HD2	1.64	0.61
1:A:788:PRO:HD3	1:A:825:MET:HE2	1.90	0.53
1:A:777:PRO:HG3	1:A:814:MET:HE2	1.91	0.53
1:B:696:LEU:HD11	1:B:755:TYR:CE2	2.43	0.52
1:B:786:ASN:ND2	1:B:786:ASN:H	1.98	0.51
1:B:786:ASN:N	1:B:786:ASN:HD22	1.92	0.51
1:A:909:GLY:O	1:A:913:MET:HG3	2.11	0.50
1:A:642:ALA:HB1	1:A:649:LEU:HD21	1.93	0.50
1:B:909:GLY:O	1:B:913:MET:HG2	2.13	0.49
1:A:863:ARG:HA	1:A:866:ARG:NH1	2.28	0.49
1:B:683:LEU:HD12	1:B:698:PRO:HA	1.94	0.48
1:B:858:ILE:HD12	1:B:869:LEU:HB2	1.95	0.48
1:A:835:SER:CB	1:A:838:HIS:HD2	2.27	0.47
1:B:894:LEU:HD12	1:B:932:MET:HE1	1.96	0.47
1:A:803:ASP:HB2	1:A:887:ASP:HB3	1.96	0.46
1:A:769:GLU:OE2	1:A:824:ARG:HD3	2.15	0.46
1:A:895:PRO:HD3	1:A:932:MET:HE3	1.98	0.46
1:A:777:PRO:HB2	1:A:811:PRO:HB3	1.98	0.46
1:A:775:ILE:O	1:A:777:PRO:HD3	2.16	0.46
1:A:894:LEU:HD12	1:A:932:MET:HE3	1.97	0.45
1:A:643:THR:O	1:A:646:GLN:O	2.34	0.45
1:A:682:VAL:HG13	1:A:695:VAL:CG1	2.45	0.45
1:B:709:ILE:HG13	1:B:710:ASN:ND2	2.31	0.45
1:B:777:PRO:HG3	1:B:814:MET:HE2	1.99	0.45
1:B:786:ASN:N	1:B:786:ASN:ND2	2.63	0.45
1:A:875:GLU:OE2	1:B:754:ARG:NH1	2.50	0.44
1:B:892:ILE:HA	1:B:892:ILE:HD13	1.82	0.44
1:A:808:THR:C	1:A:811:PRO:HD2	2.38	0.44
1:B:892:ILE:HD11	1:B:936:VAL:HG21	2.00	0.43
1:B:643:THR:O	1:B:646:GLN:O	2.36	0.43
1:B:677:ASN:HA	1:B:682:VAL:O	2.19	0.43
1:B:650:ASN:HA	1:B:691:SER:O	2.20	0.42
1:A:749:THR:O	1:A:753:GLN:HG3	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:838:HIS:O	1:B:839:GLU:C	2.58	0.42
1:A:759:ILE:O	1:A:762:GLN:HB3	2.20	0.41
1:B:903:GLU:O	1:B:906:GLU:HB2	2.20	0.41
1:B:787:GLU:HA	1:B:788:PRO:HD3	1.96	0.41
1:A:692:GLU:H	1:A:692:GLU:HG2	1.45	0.41
1:B:835:SER:HB3	1:B:868:TYR:CZ	2.56	0.41
1:A:632:TRP:CD1	1:A:663:ILE:HD11	2.57	0.40

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	294/332 (89%)	280 (95%)	11 (4%)	3 (1%)	22	63
1	B	295/332 (89%)	272 (92%)	20 (7%)	3 (1%)	22	63
All	All	589/664 (89%)	552 (94%)	31 (5%)	6 (1%)	22	63

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	841	ALA
1	A	822	LEU
1	B	822	LEU
1	A	841	ALA
1	B	839	GLU
1	A	852	SER

### 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	246/301 (82%)	229 (93%)	17 (7%)	22	54
1	B	239/301 (79%)	224 (94%)	15 (6%)	25	60
All	All	485/602 (81%)	453 (93%)	32 (7%)	24	57

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	651	ARG
1	A	671	THR
1	A	682	VAL
1	A	690	LEU
1	A	692	GLU
1	A	705	VAL
1	A	756	LEU
1	A	762	GLN
1	A	799	ILE
1	A	808	THR
1	A	822	LEU
1	A	824	ARG
1	A	851	ASP
1	A	853	THR
1	A	868	TYR
1	A	894	LEU
1	A	926	VAL
1	B	652	THR
1	B	661	SER
1	B	682	VAL
1	B	690	LEU
1	B	692	GLU
1	B	701	VAL
1	B	719	LEU
1	B	722	SER
1	B	730	GLU
1	B	750	LEU
1	B	754	ARG
1	B	786	ASN
1	B	804	TYR
1	B	868	TYR
1	B	871	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	716	ASN
1	A	718	HIS
1	A	838	HIS
1	A	859	GLN
1	A	880	ASN
1	B	659	ASN
1	B	786	ASN
1	B	820	GLN

### 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 ⓘ

There are no ligands in this entry.

### 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	298/332 (89%)	-0.14	3 (1%) 79 86	53, 64, 75, 81	0
1	B	299/332 (90%)	-0.01	3 (1%) 79 86	53, 63, 74, 86	0
All	All	597/664 (89%)	-0.08	6 (1%) 79 86	53, 64, 74, 86	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	733	ILE	3.9
1	B	652	THR	2.9
1	A	734	ARG	2.4
1	B	645	ALA	2.3
1	B	666	PRO	2.1
1	A	784	ILE	2.0

### 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 ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.