



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:18 PM GMT

PDB ID : 1KMT  
Title : Crystal structure of RhoGDI Glu(154,155)Ala mutant  
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Deposited on : 2001-12-17  
Resolution : 1.30 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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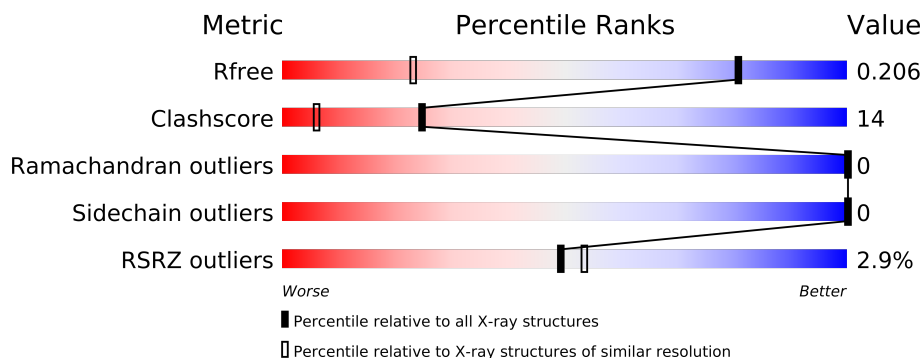
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.30 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	1025 (1.34-1.26)
Clashscore	79885	1140 (1.34-1.26)
Ramachandran outliers	78287	1093 (1.34-1.26)
Sidechain outliers	78261	1092 (1.34-1.26)
RSRZ outliers	66119	1025 (1.34-1.26)

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	
1	B	141	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4988 atoms, of which 2200 are hydrogens 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 Rho GDP-dissociation inhibitor 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	138	Total	C	H	N	O	S	78	0	0
			2198	704	1100	183	206	5			
1	B	138	Total	C	H	N	O	S	78	0	0
			2198	704	1100	183	206	5			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	GLY	-	CLONING ARTIFACT	UNP P52565
A	65	ALA	-	CLONING ARTIFACT	UNP P52565
A	66	MET	-	CLONING ARTIFACT	UNP P52565
A	154	ALA	GLU	ENGINEERED	UNP P52565
A	155	ALA	GLU	ENGINEERED	UNP P52565
B	64	GLY	-	CLONING ARTIFACT	UNP P52565
B	65	ALA	-	CLONING ARTIFACT	UNP P52565
B	66	MET	-	CLONING ARTIFACT	UNP P52565
B	154	ALA	GLU	ENGINEERED	UNP P52565
B	155	ALA	GLU	ENGINEERED	UNP P52565

- Molecule 2 is water.

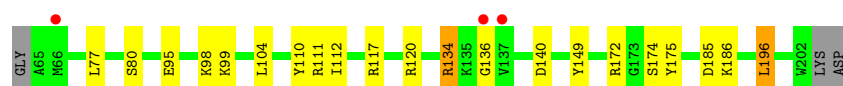
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	305	Total	O	0	0
			305	305		
2	B	287	Total	O	0	0
			287	287		

### 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.

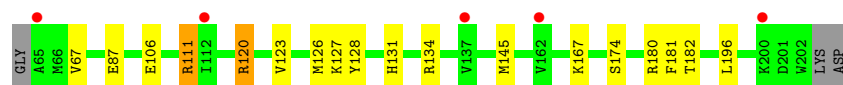
- Molecule 1: Rho GDP-dissociation inhibitor 1

Chain A: 



- Molecule 1: Rho GDP-dissociation inhibitor 1

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	34.10Å 35.92Å 67.51Å 79.19° 82.66° 76.42°	Depositor
Resolution (Å)	25.00 – 1.30 24.45 – 1.25	Depositor EDS
% Data completeness (in resolution range)	93.0 (25.00-1.30) 88.9 (24.45-1.25)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.02	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.25 (at 1.25Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.158 , 0.195 0.171 , 0.206	Depositor DCC
$R_{free}$ test set	3726 reflections (5.23%)	DCC
Wilson B-factor (Å <sup>2</sup> )	14.0	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 50.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 75030 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4988	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.90% 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	0.30	0/1123	1.28	10/1518 (0.7%)
1	B	0.32	0/1123	1.35	6/1518 (0.4%)
All	All	0.31	0/2246	1.31	16/3036 (0.5%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	ARG	NE-CZ-NH1	19.14	129.87	120.30
1	B	120	ARG	NE-CZ-NH2	-17.97	111.31	120.30
1	A	172	ARG	NE-CZ-NH2	-14.36	113.12	120.30
1	A	134	ARG	NE-CZ-NH2	-11.98	114.31	120.30
1	A	117	ARG	NE-CZ-NH2	-8.19	116.20	120.30
1	A	172	ARG	NE-CZ-NH1	8.05	124.32	120.30
1	B	134	ARG	NE-CZ-NH2	-7.61	116.49	120.30
1	A	111	ARG	CG-CD-NE	-6.98	97.13	111.80
1	A	134	ARG	NE-CZ-NH1	6.85	123.73	120.30
1	A	140	ASP	CB-CG-OD2	6.71	124.34	118.30
1	B	111	ARG	NE-CZ-NH1	-6.30	117.15	120.30
1	A	185	ASP	CB-CG-OD2	6.28	123.95	118.30
1	B	120	ARG	CD-NE-CZ	6.27	132.38	123.60
1	B	196	LEU	CA-CB-CG	6.20	129.56	115.30
1	A	117	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	A	196	LEU	CA-CB-CG	5.37	127.66	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	1098	1100	0	19	0
1	B	1098	1100	0	40	0
2	A	305	0	0	10	4
2	B	287	0	0	27	3
All	All	2788	2200	0	59	4

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 14.

All (59) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:180:ARG:HD3	2:B:447:HOH:O	1.54	1.05
1:B:123:VAL:HG12	1:B:126:MET:SD	1.97	1.04
1:B:126:MET:CE	1:B:181:PHE:CE1	2.42	1.02
1:B:182:THR:HG21	2:B:468:HOH:O	1.63	0.97
1:B:127:LYS:HD2	2:B:414:HOH:O	1.64	0.95
1:B:180:ARG:HG3	2:B:371:HOH:O	1.70	0.92
1:A:136:GLY:HA2	2:A:493:HOH:O	1.68	0.92
1:A:186:LYS:HE2	2:A:422:HOH:O	1.68	0.91
1:B:126:MET:CE	1:B:181:PHE:HE1	1.88	0.87
1:B:180:ARG:CG	2:B:371:HOH:O	2.22	0.84
1:B:87:GLU:CD	2:B:373:HOH:O	2.17	0.82
1:B:128:TYR:HA	2:B:413:HOH:O	1.81	0.81
1:B:123:VAL:CG1	1:B:126:MET:SD	2.70	0.80
1:B:131:HIS:HD2	2:B:409:HOH:O	1.64	0.79
1:B:128:TYR:HD1	2:B:413:HOH:O	1.66	0.78
1:B:87:GLU:CG	2:B:373:HOH:O	2.33	0.76
1:A:174:SER:HB3	2:A:331:HOH:O	1.85	0.75
1:B:131:HIS:CD2	2:B:409:HOH:O	2.40	0.71
1:B:127:LYS:HE2	1:B:145:MET:HG2	1.74	0.70
1:B:167:LYS:NZ	2:B:393:HOH:O	2.27	0.68
1:B:181:PHE:CD2	2:B:413:HOH:O	2.47	0.67
1:B:127:LYS:CD	2:B:414:HOH:O	2.32	0.67
1:B:128:TYR:CD1	2:B:413:HOH:O	2.42	0.67
1:B:126:MET:HE2	1:B:181:PHE:CE1	2.28	0.66
1:B:126:MET:HE1	1:B:181:PHE:CE1	2.27	0.64
1:A:136:GLY:CA	2:A:493:HOH:O	2.37	0.63
1:B:127:LYS:CE	2:B:414:HOH:O	2.48	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:174:SER:HB3	2:B:274:HOH:O	2.00	0.61
1:B:181:PHE:HA	2:B:413:HOH:O	2.02	0.58
1:B:126:MET:HE3	1:B:181:PHE:CE1	2.37	0.57
1:B:127:LYS:HD3	1:B:145:MET:HA	1.85	0.57
1:B:127:LYS:HG3	2:B:390:HOH:O	2.05	0.57
1:A:99:LYS:NZ	2:A:393:HOH:O	2.17	0.57
1:A:77:LEU:HD11	1:A:112:ILE:HD13	1.88	0.56
1:B:67:VAL:HG23	1:B:120:ARG:HD3	1.86	0.56
1:A:95:GLU:OE1	1:A:98:LYS:NZ	2.40	0.55
1:A:186:LYS:CE	2:A:422:HOH:O	2.38	0.55
1:A:149:TYR:CA	2:A:452:HOH:O	2.54	0.54
1:A:134:ARG:HD2	1:A:175:TYR:CE2	2.43	0.54
1:B:126:MET:HE1	1:B:181:PHE:HE1	1.65	0.52
1:A:77:LEU:CD1	1:A:112:ILE:CD1	2.87	0.52
1:A:149:TYR:HA	2:A:452:HOH:O	2.10	0.49
1:B:128:TYR:CA	2:B:413:HOH:O	2.52	0.49
1:B:128:TYR:CB	2:B:413:HOH:O	2.61	0.48
1:A:104:LEU:HG	1:A:196:LEU:HD12	1.94	0.48
1:B:87:GLU:CB	2:B:373:HOH:O	2.61	0.47
1:A:77:LEU:HD13	1:A:112:ILE:CD1	2.45	0.47
1:B:181:PHE:HD2	2:B:413:HOH:O	1.90	0.46
1:B:174:SER:CB	2:B:274:HOH:O	2.61	0.46
1:B:126:MET:HE2	1:B:181:PHE:HE1	1.65	0.46
1:A:77:LEU:HD12	1:A:110:TYR:CD2	2.52	0.44
1:B:180:ARG:CD	2:B:447:HOH:O	2.36	0.44
1:A:77:LEU:HD13	1:A:112:ILE:HD12	2.00	0.43
1:A:186:LYS:CD	2:A:422:HOH:O	2.65	0.42
1:B:106:GLU:HB2	1:B:167:LYS:HA	2.02	0.41
1:B:111:ARG:HH11	1:B:111:ARG:HD2	1.65	0.41
1:A:80:SER:HB3	2:A:445:HOH:O	2.20	0.41
1:B:127:LYS:CG	2:B:390:HOH:O	2.67	0.40
1:A:120:ARG:HG2	2:B:288:HOH:O	2.21	0.40

All (4) 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	Distance(Å)	Clash(Å)
2:A:467:HOH:O	2:B:265:HOH:O[1_564]	2.05	0.15
2:A:440:HOH:O	2:B:303:HOH:O[1_655]	2.12	0.08
2:A:307:HOH:O	2:B:275:HOH:O[1_655]	2.17	0.03
2:A:448:HOH:O	2:A:468:HOH:O[1_545]	2.17	0.03



## 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	136/141 (96%)	134 (98%)	2 (2%)	0	100	100
1	B	136/141 (96%)	134 (98%)	2 (2%)	0	100	100
All	All	272/282 (96%)	268 (98%)	4 (2%)	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	120/123 (98%)	120 (100%)	0	100	100
1	B	120/123 (98%)	120 (100%)	0	100	100
All	All	240/246 (98%)	240 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	69	ASN
1	B	69	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 ⓘ

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	138/141 (97%)	0.26	3 (2%) 59 64	10, 17, 27, 35	0
1	B	138/141 (97%)	0.28	5 (3%) 41 44	10, 17, 28, 38	0
All	All	276/282 (97%)	0.27	8 (2%) 49 53	10, 17, 27, 38	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	137	VAL	5.3
1	A	137	VAL	5.0
1	A	136	GLY	3.3
1	B	65	ALA	2.8
1	B	200	LYS	2.6
1	B	162	VAL	2.4
1	A	66	MET	2.2
1	B	112	ILE	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 ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.