



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:25 PM GMT

PDB ID : 3TW8  
Title : GEF domain of DENND 1B in complex with Rab GTPase Rab35  
Authors : Wu, X.D.; Kummel, D.; Reinisch, K.M.  
Deposited on : 2011-09-21  
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure 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/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

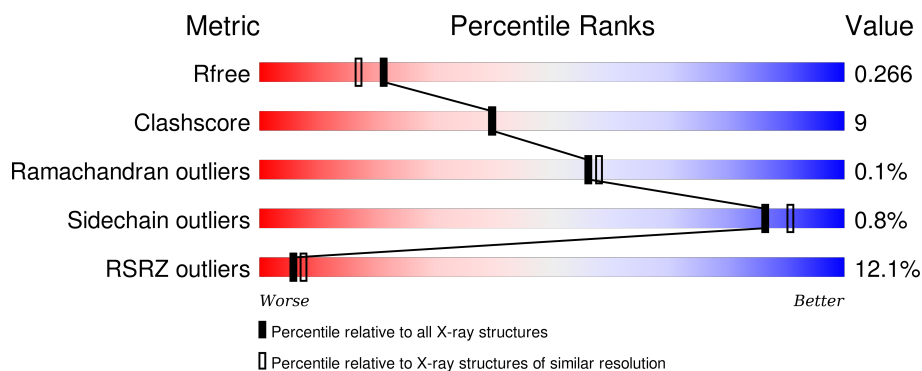
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	391	<div> <div>7%</div> <div>73% 14% • 12%</div> </div>
1	C	391	<div> <div>11%</div> <div>75% 16% • 7%</div> </div>
2	B	181	<div> <div>17%</div> <div>78% 14% • 7%</div> </div>
2	D	181	<div> <div>15%</div> <div>76% 19% • •</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8573 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DENN domain-containing protein 1B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	87	0	0
			2751	1772	458	504	17			
1	C	362	Total	C	N	O	S	115	0	0
			2893	1863	480	533	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	EXPRESSION TAG	UNP Q6P3S1-2
C	1	SER	-	EXPRESSION TAG	UNP Q6P3S1-2

- Molecule 2 is a protein called Ras-related protein Rab-35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	169	Total	C	N	O	S	60	0	0
			1357	858	232	262	5			
2	D	173	Total	C	N	O	S	75	0	0
			1385	876	236	268	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	EXPRESSION TAG	UNP Q15286
D	0	SER	-	EXPRESSION TAG	UNP Q15286

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	80	Total	O	0	0
			80	80		
3	B	18	Total	O	0	0
			18	18		

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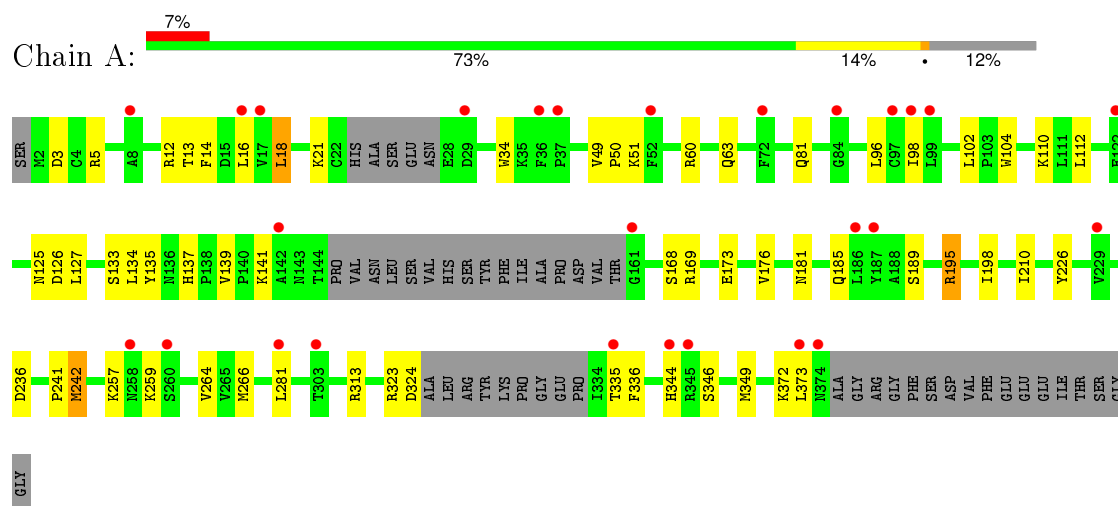
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	64	Total	O	0	0
			64	64		
3	D	25	Total	O	0	0
			25	25		

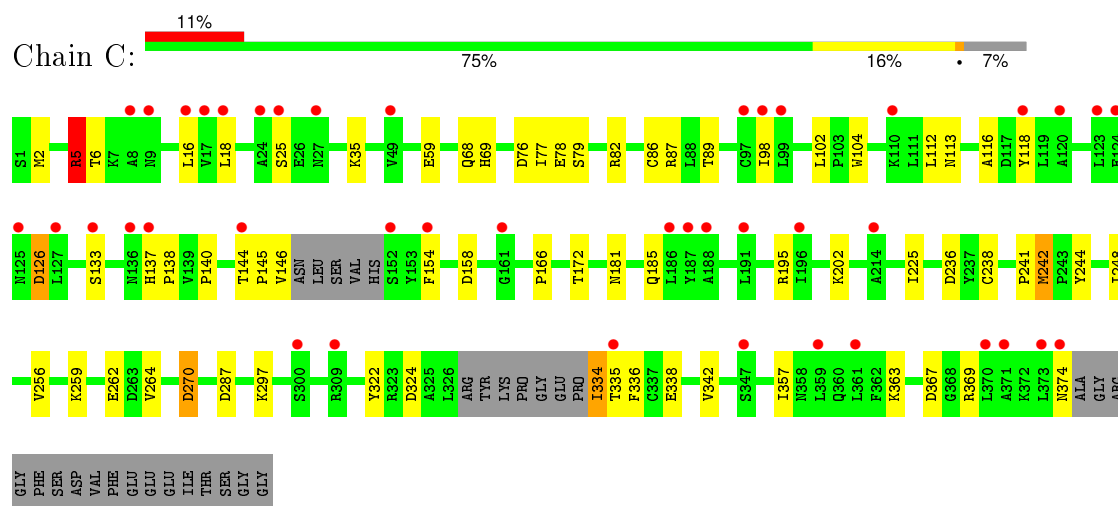
### 3 Residue-property plots

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.

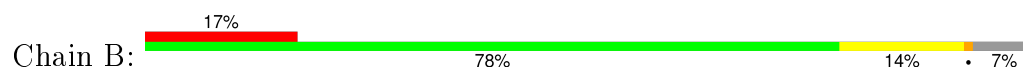
- Molecule 1: DENN domain-containing protein 1B

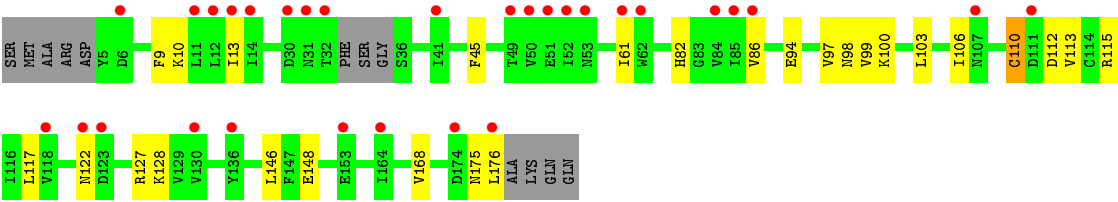


- Molecule 1: DENN domain-containing protein 1B

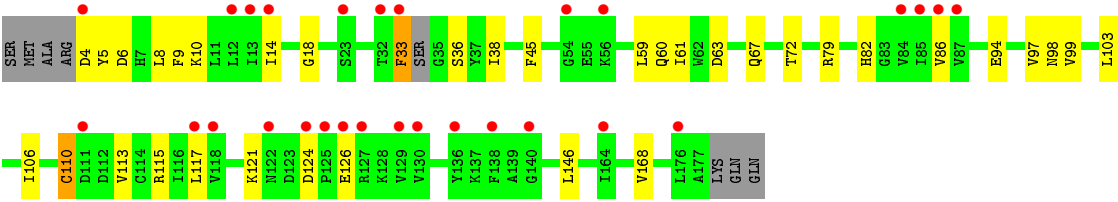
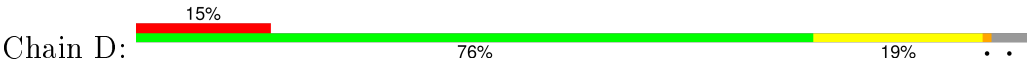


- Molecule 2: Ras-related protein Rab-35





• Molecule 2: Ras-related protein Rab-35



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.59Å 56.56Å 175.31Å 90.00° 95.63° 90.00°	Depositor
Resolution (Å)	24.89 – 2.10 24.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (24.89-2.10) 91.2 (24.89-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.57 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.227 , 0.269 0.226 , 0.266	Depositor DCC
$R_{free}$ test set	4502 reflections (7.56%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.8	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 46.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 64040 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8573	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.97	3/2812 (0.1%)	1.00	13/3813 (0.3%)
1	C	1.00	8/2960 (0.3%)	0.99	12/4020 (0.3%)
2	B	0.79	2/1377 (0.1%)	0.82	3/1858 (0.2%)
2	D	0.92	3/1406 (0.2%)	0.89	6/1897 (0.3%)
All	All	0.94	16/8555 (0.2%)	0.95	34/11588 (0.3%)

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	195	ARG	CZ-NH1	-11.05	1.18	1.33
1	A	195	ARG	CZ-NH1	-9.84	1.20	1.33
1	C	195	ARG	CZ-NH2	-9.14	1.21	1.33
1	A	195	ARG	CZ-NH2	-8.98	1.21	1.33
1	C	238	CYS	CB-SG	-8.88	1.67	1.82

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	195	ARG	NE-CZ-NH2	20.52	130.56	120.30
1	A	195	ARG	NE-CZ-NH1	15.61	128.11	120.30
1	C	195	ARG	NH1-CZ-NH2	-14.04	103.96	119.40
1	A	195	ARG	NE-CZ-NH2	13.68	127.14	120.30
1	A	195	ARG	NH1-CZ-NH2	-13.48	104.58	119.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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 hydrogen



atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2751	0	2760	49	1
1	C	2893	0	2896	47	2
2	B	1357	0	1348	19	0
2	D	1385	0	1369	28	2
3	A	80	0	0	3	0
3	B	18	0	0	1	0
3	C	64	0	0	7	0
3	D	25	0	0	1	0
All	All	8573	0	8373	141	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

The worst 5 of 141 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:LYS:NZ	1:A:266:MET:HE1	1.80	0.96
1:A:257:LYS:NZ	1:A:266:MET:CE	2.39	0.85
1:A:257:LYS:HZ1	1:A:266:MET:CE	1.91	0.83
1:C:242:MET:H	1:C:242:MET:HE3	1.47	0.77
1:C:262:GLU:O	1:C:264:VAL:HG23	1.85	0.76

All (3) 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ASN:CG	1:A:181:ASN:ND2[1_545]	2.15	0.05
1:C:287:ASP:OD2	2:D:33:PHE:CE1[1_565]	2.17	0.03
1:C:287:ASP:OD2	2:D:33:PHE:CZ[1_565]	2.18	0.02

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	335/391 (86%)	322 (96%)	12 (4%)	1 (0%)	46	45
1	C	356/391 (91%)	345 (97%)	11 (3%)	0	100	100
2	B	165/181 (91%)	158 (96%)	7 (4%)	0	100	100
2	D	169/181 (93%)	164 (97%)	5 (3%)	0	100	100
All	All	1025/1144 (90%)	989 (96%)	35 (3%)	1 (0%)	56	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	373	LEU

### 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	312/351 (89%)	308 (99%)	4 (1%)	76	82
1	C	328/351 (93%)	324 (99%)	4 (1%)	78	84
2	B	151/160 (94%)	151 (100%)	0	100	100
2	D	153/160 (96%)	153 (100%)	0	100	100
All	All	944/1022 (92%)	936 (99%)	8 (1%)	86	91

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	242	MET
1	C	334	ILE
1	C	242	MET
1	A	110	LYS
1	C	5	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such

sidechains are listed below:

Mol	Chain	Res	Type
2	B	122	ASN
1	C	69	HIS
2	D	60	GLN
2	B	82	HIS
1	C	374	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	343/391 (87%)	0.47	27 (7%) 15 21	26, 45, 83, 104	22 (6%)
1	C	362/391 (92%)	0.61	42 (11%) 6 9	26, 50, 84, 120	33 (9%)
2	B	169/181 (93%)	1.05	30 (17%) 2 2	36, 62, 86, 99	16 (9%)
2	D	173/181 (95%)	0.68	28 (16%) 3 3	31, 52, 80, 98	19 (10%)
All	All	1047/1144 (91%)	0.65	127 (12%) 6 7	26, 51, 85, 120	90 (8%)

The worst 5 of 127 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	31	ASN	5.5
2	B	52	ILE	5.4
2	D	176	LEU	5.1
1	C	8	ALA	5.0
2	B	85	ILE	4.9

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.