



# wwPDB X-ray Structure Validation Summary Report

Mar 1, 2014 – 03:48 AM GMT

PDB ID : 1L6O  
Title : XENOPUS DISHEVELLED PDZ DOMAIN  
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Deposited on : 2002-03-11  
Resolution : 2.20 Å(reported)

This is a wwPDB 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/ValidationPDFNotes.html>

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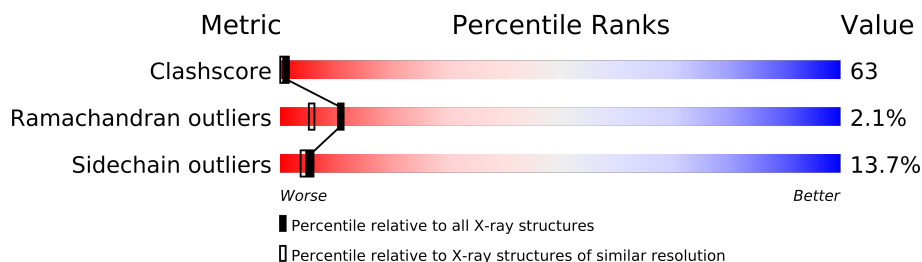
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
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 2.20 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))
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	95	
1	B	95	
1	C	95	
2	D	8	
2	E	8	
2	F	8	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2328 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 Segment polarity protein dishevelled homolog DVL-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	95	Total	C	N	O	Se	0	0	0
			724	453	128	138	5			
1	B	93	Total	C	N	O	Se	0	0	0
			703	441	122	135	5			
1	C	92	Total	C	N	O	Se	0	0	0
			693	435	119	134	5			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	MSE	-	SEE REMARK 999	UNP P51142
A	259	MSE	MET	MODIFIED RESIDUE	UNP P51142
A	287	MSE	MET	MODIFIED RESIDUE	UNP P51142
A	303	MSE	MET	MODIFIED RESIDUE	UNP P51142
A	315	MSE	MET	MODIFIED RESIDUE	UNP P51142
A	341	LEU	-	EXPRESSION TAG	UNP P51142
A	342	GLU	-	EXPRESSION TAG	UNP P51142
A	343	HIS	-	EXPRESSION TAG	UNP P51142
A	344	HIS	-	EXPRESSION TAG	UNP P51142
A	345	HIS	-	EXPRESSION TAG	UNP P51142
B	251	MSE	-	SEE REMARK 999	UNP P51142
B	259	MSE	MET	MODIFIED RESIDUE	UNP P51142
B	287	MSE	MET	MODIFIED RESIDUE	UNP P51142
B	303	MSE	MET	MODIFIED RESIDUE	UNP P51142
B	315	MSE	MET	MODIFIED RESIDUE	UNP P51142
B	341	LEU	-	EXPRESSION TAG	UNP P51142
B	342	GLU	-	EXPRESSION TAG	UNP P51142
B	343	HIS	-	EXPRESSION TAG	UNP P51142
B	344	HIS	-	EXPRESSION TAG	UNP P51142
B	345	HIS	-	EXPRESSION TAG	UNP P51142
C	251	MSE	-	SEE REMARK 999	UNP P51142
C	259	MSE	MET	MODIFIED RESIDUE	UNP P51142
C	287	MSE	MET	MODIFIED RESIDUE	UNP P51142

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Chain	Residue	Modelled	Actual	Comment	Reference
C	303	MSE	MET	MODIFIED RESIDUE	UNP P51142
C	315	MSE	MET	MODIFIED RESIDUE	UNP P51142
C	341	LEU	-	EXPRESSION TAG	UNP P51142
C	342	GLU	-	EXPRESSION TAG	UNP P51142
C	343	HIS	-	EXPRESSION TAG	UNP P51142
C	344	HIS	-	EXPRESSION TAG	UNP P51142
C	345	HIS	-	EXPRESSION TAG	UNP P51142

- Molecule 2 is a protein called Dapper 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	8	Total 61	C 39	N 9	O 12	S 1	0	0	0
2	E	8	Total 61	C 39	N 9	O 12	S 1	0	0	0
2	F	8	Total 61	C 39	N 9	O 12	S 1	0	0	0

- Molecule 3 is water.

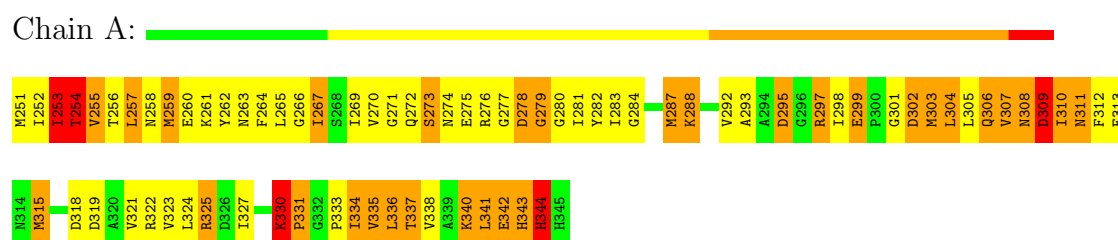
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total 16	O 16	0	0
3	B	3	Total 3	O 3	0	0
3	C	4	Total 4	O 4	0	0
3	E	1	Total 1	O 1	0	0
3	F	1	Total 1	O 1	0	0

### 3 Residue-property plots

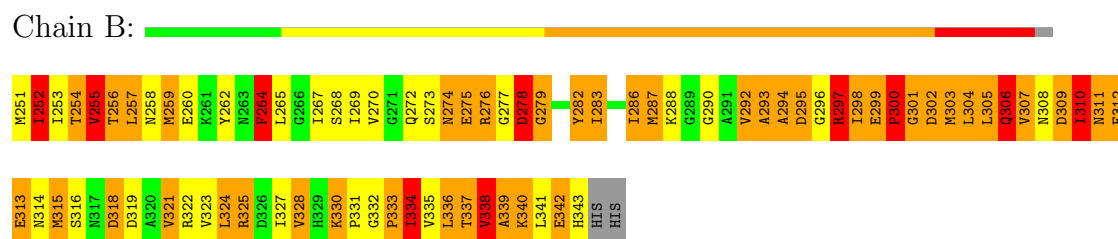
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

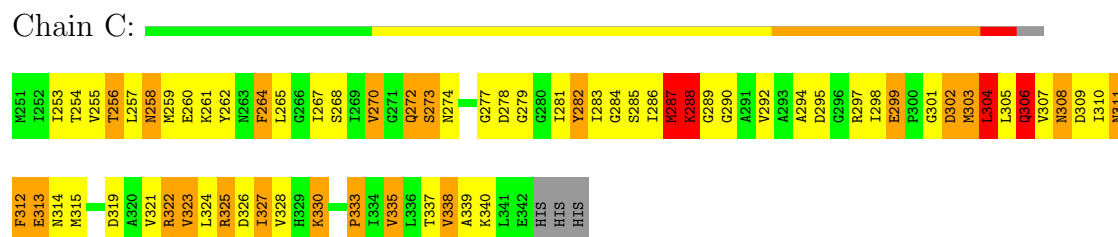
- Molecule 1: Segment polarity protein dishevelled homolog DVL-2



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- Molecule 1: Segment polarity protein dishevelled homolog DVL-2



- Molecule 2: Dapper 1



- Molecule 2: Dapper 1





- Molecule 2: Dapper 1

Chain F: 



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.73Å 84.73Å 123.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.23 – 2.20	Depositor
% Data completeness (in resolution range)	96.0 (54.23-2.20)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.277 , 0.323	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2328	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	64.0	wwPDB-VP

## 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	3.02	72/730 (9.9%)	2.33	38/977 (3.9%)
1	B	2.90	65/707 (9.2%)	2.63	50/947 (5.3%)
1	C	2.68	47/696 (6.8%)	2.18	31/932 (3.3%)
2	D	3.50	8/60 (13.3%)	2.27	3/78 (3.8%)
2	E	3.52	6/60 (10.0%)	2.95	8/78 (10.3%)
2	F	3.94	10/60 (16.7%)	2.65	4/78 (5.1%)
All	All	2.94	208/2313 (9.0%)	2.41	134/3090 (4.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3
2	D	0	1
2	F	0	1
All	All	0	5

The worst 5 of 208 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	338	VAL	CB-CG2	-17.09	1.17	1.52
1	A	303	MSE	SE-CE	-13.79	1.14	1.95
2	F	6	THR	CB-CG2	-12.63	1.10	1.52
1	A	284	GLY	C-O	-12.06	1.04	1.23
1	B	315	MSE	SE-CE	-12.04	1.24	1.95

The worst 5 of 134 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	312	PHE	CB-CG-CD1	14.96	131.27	120.80
1	B	312	PHE	CB-CG-CD2	-14.77	110.46	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	288	LYS	C-N-CA	-13.85	93.22	122.30
1	B	318	ASP	CB-CG-OD1	-13.68	105.99	118.30
1	B	324	LEU	CB-CG-CD2	-13.46	88.11	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	252	ILE	Mainchain
1	B	300	PRO	Peptide
1	B	310	ILE	Mainchain
2	D	6	THR	Mainchain
2	F	7	THR	Mainchain

## 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	724	0	727	97	0
1	B	703	0	713	89	0
1	C	693	0	705	76	0
2	D	61	0	74	18	0
2	E	61	0	74	12	0
2	F	61	0	74	19	0
3	A	16	0	0	17	0
3	B	3	0	0	2	0
3	C	4	0	0	1	0
3	E	1	0	0	0	0
3	F	1	0	0	1	0
All	All	2328	0	2367	296	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 63.

The worst 5 of 296 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:310:ILE:CG2	1:A:310:ILE:CB	1.74	1.65
1:B:324:LEU:CD1	1:B:324:LEU:CG	1.77	1.60
2:E:6:THR:CA	2:E:6:THR:CB	1.76	1.59
2:D:7:THR:CB	2:D:7:THR:CA	1.74	1.58
1:C:325:ARG:CG	1:C:325:ARG:CD	1.76	1.58

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	93/95 (98%)	87 (94%)	4 (4%)	2 (2%)	10	5
1	B	91/95 (96%)	75 (82%)	12 (13%)	4 (4%)	4	1
1	C	90/95 (95%)	82 (91%)	8 (9%)	0	100	100
2	D	6/8 (75%)	5 (83%)	1 (17%)	0	100	100
2	E	6/8 (75%)	6 (100%)	0	0	100	100
2	F	6/8 (75%)	6 (100%)	0	0	100	100
All	All	292/309 (94%)	261 (89%)	25 (9%)	6 (2%)	11	5

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	344	HIS
1	B	301	GLY
1	A	309	ASP
1	B	278	ASP
1	B	275	GLU

### 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	79/74 (107%)	70 (89%)	9 (11%)	8	7
1	B	77/74 (104%)	65 (84%)	12 (16%)	4	3
1	C	76/74 (103%)	67 (88%)	9 (12%)	8	6
2	D	8/8 (100%)	8 (100%)	0	100	100
2	E	8/8 (100%)	6 (75%)	2 (25%)	1	0
2	F	8/8 (100%)	5 (62%)	3 (38%)	0	0
All	All	256/246 (104%)	221 (86%)	35 (14%)	5	4

5 of 35 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	287	MSE
1	B	337	THR
2	F	1	SER
1	B	297	ARG
1	B	306	GLN

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

Mol	Chain	Res	Type
1	B	308	ASN
1	B	311	ASN
1	C	272	GLN
1	B	272	GLN
1	C	258	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 ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.