



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 11:05 AM UTC

PDB ID : 1L6O / pdb_00001l6o
Title : XENOPUS DISHEVELLED PDZ DOMAIN
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Khlebtsova, N.; Fox, E.P.; Earnest, T.; Moon, R.T.
Deposited on : 2002-03-11
Resolution : 2.20 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

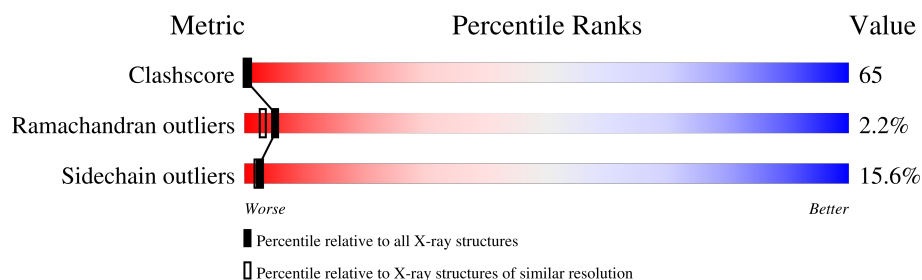
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	95	8% 45% 37% 9%
1	B	95	7% 29% 46% 15% .
1	C	95	9% 48% 34% 5% .
2	D	8	12% 25% 50% 12%
2	E	8	12% 25% 38% 25%
2	F	8	12% 25% 25% 38%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2328 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 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	C	N	O	S	0	0	0
			61	39	9	12	1			
2	E	8	Total	C	N	O	S	0	0	0
			61	39	9	12	1			
2	F	8	Total	C	N	O	S	0	0	0
			61	39	9	12	1			

- Molecule 3 is water.

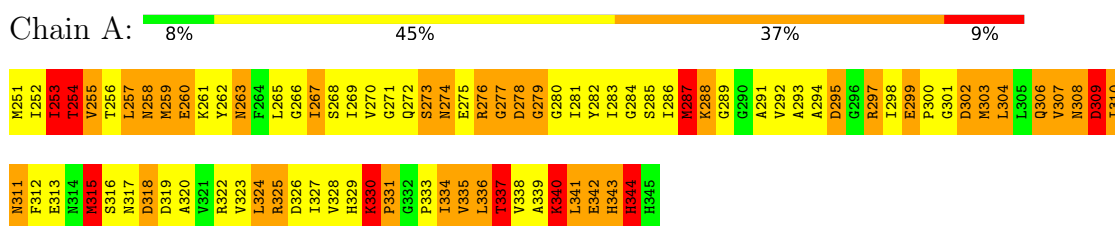
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	16	Total	O	0	0
			16	16		
3	B	3	Total	O	0	0
			3	3		
3	C	4	Total	O	0	0
			4	4		
3	E	1	Total	O	0	0
			1	1		
3	F	1	Total	O	0	0
			1	1		

3 Residue-property plots

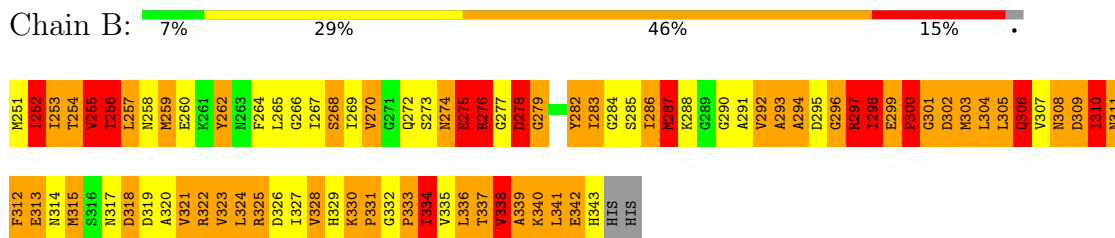
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

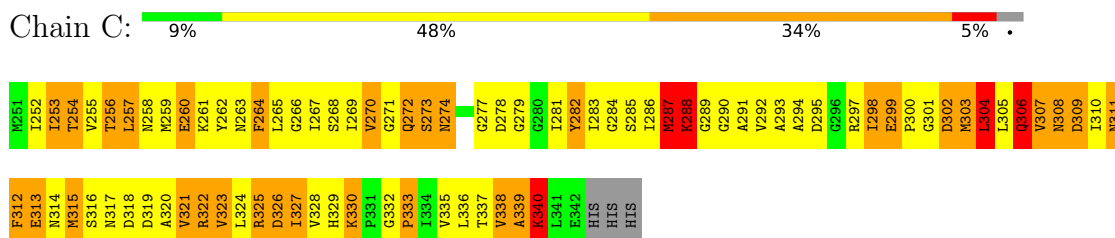
- Molecule 1: Segment polarity protein dishevelled homolog DVL-2



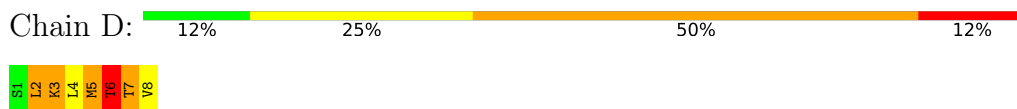
- Molecule 1: Segment polarity protein dishevelled homolog DVL-2



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



- Molecule 2: Dapper 1



- Molecule 2: Dapper 1





● Molecule 2: Dapper 1



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	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 5.0	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 ²)	64.0	wwPDB-VP

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	4.20	146/730 (20.0%)	2.81	74/977 (7.6%)
1	B	3.95	110/707 (15.6%)	3.10	108/947 (11.4%)
1	C	3.61	101/696 (14.5%)	2.87	60/932 (6.4%)
2	D	5.06	11/60 (18.3%)	2.88	8/78 (10.3%)
2	E	5.08	9/60 (15.0%)	2.98	7/78 (9.0%)
2	F	5.32	12/60 (20.0%)	3.18	8/78 (10.3%)
All	All	4.04	389/2313 (16.8%)	2.93	265/3090 (8.6%)

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 389 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	303	MSE	SE-CE	-27.11	1.14	1.95
2	D	5	MET	SD-CE	-24.34	1.18	1.79
1	B	315	MSE	SE-CE	-23.68	1.24	1.95
2	E	5	MET	SD-CE	-22.39	1.23	1.79
1	C	286	ILE	C-O	-20.29	1.02	1.24

The worst 5 of 265 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	288	LYS	CA-C-N	-19.17	93.22	123.31
1	C	288	LYS	C-N-CA	-19.17	93.22	123.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	288	LYS	N-CA-CB	17.50	134.09	110.29
1	B	334	ILE	CB-CA-C	-11.76	92.01	111.29
1	C	288	LYS	O-C-N	-11.50	109.11	122.68

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 Too-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 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	724	0	727	99	0
1	B	703	0	713	93	0
1	C	693	0	705	79	0
2	D	61	0	74	18	0
2	E	61	0	74	12	0
2	F	61	0	74	17	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	303	0

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 65.

The worst 5 of 303 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:310:ILE:CG2	1:A:310:ILE:CB	1.74	1.65
1:B:303:MSE:C	1:B:303:MSE:CA	1.74	1.60
1:B:324:LEU:CG	1:B:324:LEU:CD1	1.77	1.60
2:E:6:THR:CB	2:E:6:THR:CA	1.76	1.59
1:C:325:ARG:CD	1:C:325:ARG:CG	1.76	1.58

There are no symmetry-related clashes.

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	81/95 (85%)	75 (93%)	4 (5%)	2 (2%)	4	2
1	B	35/95 (37%)	28 (80%)	4 (11%)	3 (9%)	0	0
1	C	89/95 (94%)	81 (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	223/309 (72%)	201 (90%)	17 (8%)	5 (2%)	5	3

All (5) 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 [i](#)

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%)	67 (85%)	12 (15%)	3	2
1	B	77/74 (104%)	63 (82%)	14 (18%)	2	1
1	C	76/74 (103%)	67 (88%)	9 (12%)	5	5
2	D	8/8 (100%)	8 (100%)	0	100	100
2	E	8/8 (100%)	6 (75%)	2 (25%)	0	0
2	F	8/8 (100%)	5 (62%)	3 (38%)	0	0
All	All	256/246 (104%)	216 (84%)	40 (16%)	2	2

5 of 40 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	287	MSE
2	E	3	LYS
1	C	288	LYS
1	C	321	VAL
2	F	1	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	258	ASN
1	C	272	GLN
1	C	308	ASN
1	C	306	GLN
1	B	263	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 oligosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	288:LYS	C	289:GLY	N	1.18

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.