



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 2, 2016 – 12:04 PM EDT

PDB ID : 5GLF  
Title : Structural insights into the interaction of p97 N-terminal domain and SHP motif in Derlin-1 rhomboid pseudoprotease  
Authors : Lim, J.J.; Lee, Y.; Yoon, S.Y.; Ly, T.T.; Kang, J.Y.; Youn, H.-S.; An, J.Y.; Lee, J.-G.; Park, K.R.; Kim, T.G.; Yang, J.K.; Jun, Y.; Eom, S.H.  
Deposited on : 2016-07-11  
Resolution : 2.25 Å(reported)

This is a Full wwPDB X-ray Structure 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/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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028320
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)	:	rb-20028320

**i**

## X-RAY DIFFRACTION



A.

Ramachandran outliers

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.


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Mol	Chain	Length	Quality of chain
2	F	12	 67% 17% 17%
2	H	12	 17% 83% 17%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6171 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 Transitional endoplasmic reticulum ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1355	855	240	252	8			
1	C	171	Total	C	N	O	S	0	0	0
			1366	861	244	253	8			
1	E	169	Total	C	N	O	S	0	0	0
			1348	850	239	251	8			
1	G	168	Total	C	N	O	S	0	0	0
			1335	844	235	248	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	-	expression tag	UNP P55072
A	17	ALA	-	expression tag	UNP P55072
A	18	MET	-	expression tag	UNP P55072
A	19	GLY	-	expression tag	UNP P55072
A	20	SER	-	expression tag	UNP P55072
C	16	GLY	-	expression tag	UNP P55072
C	17	ALA	-	expression tag	UNP P55072
C	18	MET	-	expression tag	UNP P55072
C	19	GLY	-	expression tag	UNP P55072
C	20	SER	-	expression tag	UNP P55072
E	16	GLY	-	expression tag	UNP P55072
E	17	ALA	-	expression tag	UNP P55072
E	18	MET	-	expression tag	UNP P55072
E	19	GLY	-	expression tag	UNP P55072
E	20	SER	-	expression tag	UNP P55072
G	16	GLY	-	expression tag	UNP P55072
G	17	ALA	-	expression tag	UNP P55072
G	18	MET	-	expression tag	UNP P55072
G	19	GLY	-	expression tag	UNP P55072
G	20	SER	-	expression tag	UNP P55072

- Molecule 2 is a protein called Derlin-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	12	Total	C	N	O	0	0	0
			102	63	23	16			
2	D	11	Total	C	N	O	0	0	0
			91	57	19	15			
2	F	10	Total	C	N	O	0	0	0
			83	53	18	12			
2	H	10	Total	C	N	O	0	0	0
			83	53	18	12			

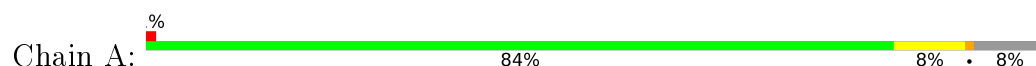
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	117	Total	O	0	0
			117	117		
3	B	10	Total	O	0	0
			10	10		
3	C	115	Total	O	0	0
			115	115		
3	D	11	Total	O	0	0
			11	11		
3	E	71	Total	O	0	0
			71	71		
3	F	6	Total	O	0	0
			6	6		
3	G	76	Total	O	0	0
			76	76		
3	H	2	Total	O	0	0
			2	2		

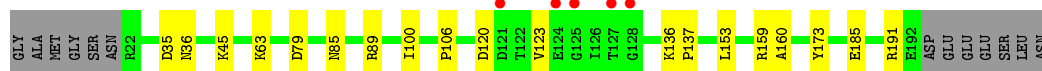
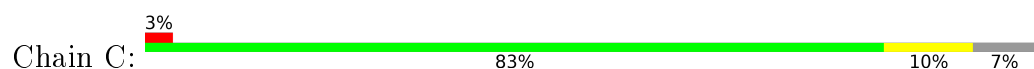
### 3 Residue-property plots [i](#)

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.

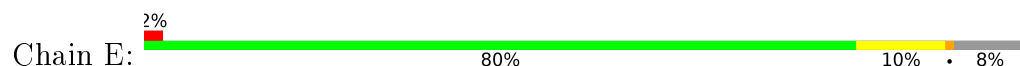
- Molecule 1: Transitional endoplasmic reticulum ATPase



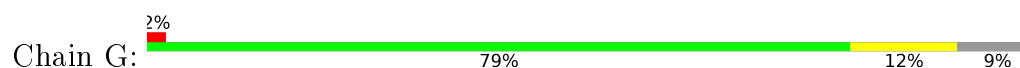
- Molecule 1: Transitional endoplasmic reticulum ATPase



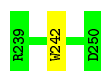
- Molecule 1: Transitional endoplasmic reticulum ATPase



- Molecule 1: Transitional endoplasmic reticulum ATPase



- Molecule 2: Derlin-1



- Molecule 2: Derlin-1

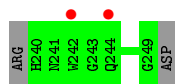
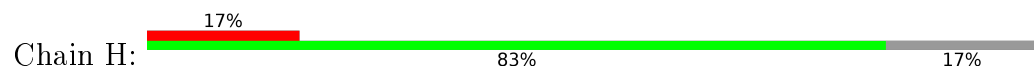




- Molecule 2: Derlin-1



- Molecule 2: Derlin-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	43.80 Å 74.40 Å 222.11 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.97 – 2.25 42.97 – 2.25	Depositor EDS
% Data completeness (in resolution range)	96.6 (42.97-2.25) 96.7 (42.97-2.25)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.11 (at 2.24 Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.189 , 0.230 0.196 , 0.236	Depositor DCC
$R_{free}$ test set	1716 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	30.7	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 46.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6171	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 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	0.80	0/1377	0.92	2/1861 (0.1%)
1	C	0.83	0/1388	0.97	2/1876 (0.1%)
1	E	0.70	0/1369	0.99	5/1850 (0.3%)
1	G	0.69	0/1357	0.93	1/1835 (0.1%)
2	B	0.87	0/105	0.87	0/139
2	D	0.68	0/94	0.60	0/125
2	F	0.76	0/86	0.72	0/114
2	H	0.68	0/86	0.67	0/114
All	All	0.76	0/5862	0.94	10/7914 (0.1%)

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	A	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	GLY	N-CA-C	7.09	130.81	113.10
1	E	157	GLY	N-CA-C	-7.01	95.56	113.10
1	C	89	ARG	NE-CZ-NH2	6.78	123.69	120.30
1	E	191	ARG	NE-CZ-NH1	-5.63	117.48	120.30
1	G	89	ARG	NE-CZ-NH2	5.59	123.09	120.30
1	C	173	TYR	CB-CA-C	-5.47	99.47	110.40
1	A	25	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	E	120	ASP	CB-CG-OD1	5.37	123.13	118.30
1	E	191	ARG	CG-CD-NE	-5.27	100.73	111.80
1	E	74	ASP	CB-CG-OD1	5.15	122.94	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	157	GLY	Peptide

## 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	1355	0	1391	9	0
1	C	1366	0	1403	13	0
1	E	1348	0	1383	15	0
1	G	1335	0	1372	21	0
2	B	102	0	89	1	0
2	D	91	0	76	0	0
2	F	83	0	72	1	0
2	H	83	0	72	0	0
3	A	117	0	0	1	0
3	B	10	0	0	0	0
3	C	115	0	0	3	0
3	D	11	0	0	1	0
3	E	71	0	0	1	0
3	F	6	0	0	0	0
3	G	76	0	0	2	0
3	H	2	0	0	0	0
All	All	6171	0	5858	56	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 5.

All (56) 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:G:123:VAL:HA	1:G:126:ILE:HD12	1.35	1.04
1:C:120:ASP:O	1:C:123:VAL:HG22	1.73	0.88
1:E:24:ASN:ND2	1:E:49:LEU:HD21	1.98	0.79
1:G:123:VAL:HA	1:G:126:ILE:CD1	2.12	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:123:VAL:CA	1:G:126:ILE:HD12	2.14	0.76
1:G:129:ASN:HB3	1:G:132:GLU:HB2	1.69	0.74
1:C:106:PRO:HD2	3:C:233:HOH:O	1.87	0.73
1:E:153:LEU:HD11	1:E:160:ALA:HB1	1.68	0.73
1:G:120:ASP:CG	1:G:190:LYS:HB3	2.14	0.67
1:E:121:ASP:OD2	1:E:191:ARG:HG3	1.95	0.67
1:C:36:ASN:HB3	3:C:224:HOH:O	1.96	0.64
1:E:158:MET:O	1:E:158:MET:HG3	1.97	0.63
1:A:24:ASN:CG	1:A:49:LEU:HD21	2.20	0.62
1:E:120:ASP:OD2	1:E:190:LYS:HD2	2.01	0.61
1:C:45:LYS:HE3	1:C:79:ASP:O	2.02	0.59
1:G:95:ARG:HG3	3:G:273:HOH:O	2.04	0.58
1:C:36:ASN:HA	1:C:85:ASN:HD21	1.69	0.56
1:C:35:ASP:OD2	1:E:64:ARG:NH2	2.38	0.56
1:G:189:ILE:O	1:G:190:LYS:HB2	2.04	0.56
1:A:158:MET:HG3	1:A:158:MET:O	2.06	0.56
1:G:120:ASP:OD2	1:G:190:LYS:HB3	2.05	0.56
1:G:36:ASN:HA	1:G:85:ASN:HD21	1.70	0.55
1:A:136:LYS:HB3	1:A:137:PRO:HD3	1.90	0.54
1:E:70:ILE:HG12	3:E:235:HOH:O	2.07	0.54
1:E:36:ASN:HA	1:E:85:ASN:HD21	1.72	0.54
1:E:143:TYR:CZ	1:G:63:LYS:HE2	2.42	0.53
1:A:107:ASP:HB2	3:A:266:HOH:O	2.09	0.53
1:C:120:ASP:OD1	1:C:191:ARG:NH1	2.36	0.52
1:C:136:LYS:HB3	1:C:137:PRO:HD3	1.93	0.51
1:G:120:ASP:OD1	1:G:190:LYS:HB3	2.10	0.50
2:F:247:ARG:NH1	2:F:249:GLY:HA2	2.27	0.50
1:C:159:ARG:HD2	3:C:216:HOH:O	2.12	0.49
1:E:143:TYR:CE2	1:G:63:LYS:HE2	2.47	0.49
1:C:185:GLU:HB2	3:D:303:HOH:O	2.13	0.48
1:E:136:LYS:HB3	1:E:137:PRO:HD3	1.95	0.48
1:G:136:LYS:HB3	1:G:137:PRO:HD3	1.96	0.48
1:A:34:GLU:O	1:A:34:GLU:HG3	2.13	0.47
1:C:153:LEU:HD11	1:C:160:ALA:HB1	1.96	0.47
1:G:78:SER:HB3	1:G:80:GLU:OE1	2.15	0.47
1:G:103:GLN:HB3	1:G:104:PRO:HD2	1.97	0.46
1:C:63:LYS:HB3	1:C:63:LYS:HE2	1.88	0.44
1:G:189:ILE:O	1:G:190:LYS:CB	2.65	0.44
1:A:167:GLU:HG2	2:B:242:TRP:CZ2	2.54	0.43
1:G:103:GLN:HG3	3:G:231:HOH:O	2.19	0.43
1:E:64:ARG:HG2	1:E:64:ARG:HH11	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:120:ASP:OD2	1:E:190:LYS:CD	2.67	0.42
1:G:100:ILE:C	1:G:100:ILE:HD12	2.40	0.42
1:G:123:VAL:HG12	1:G:126:ILE:HD12	2.01	0.42
1:G:120:ASP:OD1	1:G:121:ASP:N	2.53	0.42
1:G:65:ARG:HG3	1:G:93:ARG:HG2	2.01	0.41
1:A:120:ASP:OD1	1:A:121:ASP:N	2.54	0.41
1:E:100:ILE:HD12	1:E:100:ILE:C	2.41	0.41
1:A:100:ILE:HD12	1:A:100:ILE:C	2.40	0.41
1:A:134:TYR:CE1	1:A:156:GLY:HA3	2.56	0.40
1:C:100:ILE:HD12	1:C:100:ILE:C	2.42	0.40
1:E:65:ARG:HG3	1:E:93:ARG:HG2	2.03	0.40

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	168/184 (91%)	165 (98%)	3 (2%)	0	100	100
1	C	169/184 (92%)	164 (97%)	5 (3%)	0	100	100
1	E	167/184 (91%)	164 (98%)	3 (2%)	0	100	100
1	G	166/184 (90%)	164 (99%)	2 (1%)	0	100	100
2	B	10/12 (83%)	10 (100%)	0	0	100	100
2	D	9/12 (75%)	9 (100%)	0	0	100	100
2	F	8/12 (67%)	8 (100%)	0	0	100	100
2	H	8/12 (67%)	8 (100%)	0	0	100	100
All	All	705/784 (90%)	692 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	155/166 (93%)	155 (100%)	0	100	100
1	C	156/166 (94%)	156 (100%)	0	100	100
1	E	154/166 (93%)	154 (100%)	0	100	100
1	G	153/166 (92%)	153 (100%)	0	100	100
2	B	9/9 (100%)	9 (100%)	0	100	100
2	D	8/9 (89%)	8 (100%)	0	100	100
2	F	7/9 (78%)	7 (100%)	0	100	100
2	H	7/9 (78%)	7 (100%)	0	100	100
All	All	649/700 (93%)	649 (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 (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	24	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 ⓘ

### 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	170/184 (92%)	-0.40	2 (1%) 81 83	15, 24, 52, 86	0
1	C	171/184 (92%)	-0.39	5 (2%) 55 60	19, 27, 66, 85	0
1	E	169/184 (91%)	-0.16	3 (1%) 71 75	28, 39, 66, 89	0
1	G	168/184 (91%)	-0.20	3 (1%) 71 75	22, 41, 78, 93	0
2	B	12/12 (100%)	-0.28	0 100 100	24, 33, 45, 66	0
2	D	11/12 (91%)	-0.48	0 100 100	23, 30, 45, 47	0
2	F	10/12 (83%)	-0.06	0 100 100	39, 47, 58, 59	0
2	H	10/12 (83%)	1.11	2 (20%) 1 1	42, 58, 67, 69	0
All	All	721/784 (91%)	-0.27	15 (2%) 67 71	15, 35, 67, 93	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	127	THR	6.4
1	E	156	GLY	6.2
1	E	158	MET	5.4
1	E	34	GLU	3.5
1	A	34	GLU	3.2
1	G	127	THR	3.0
1	A	158	MET	2.9
1	C	128	GLY	2.8
1	C	121	ASP	2.8
1	G	157	GLY	2.4
1	C	124	GLU	2.4
1	C	125	GLY	2.2
2	H	244	GLN	2.2
2	H	242	TRP	2.1
1	G	123	VAL	2.0

## 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.