



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3GRL / pdb_00003grl
Title : Crystal Structure of the Monomer of the p115 Tether Globular Head Domain
Authors : An, Y.; Elsliger, M.A.; Wilson, I.A.
Deposited on : 2009-03-25
Resolution : 2.00 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

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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

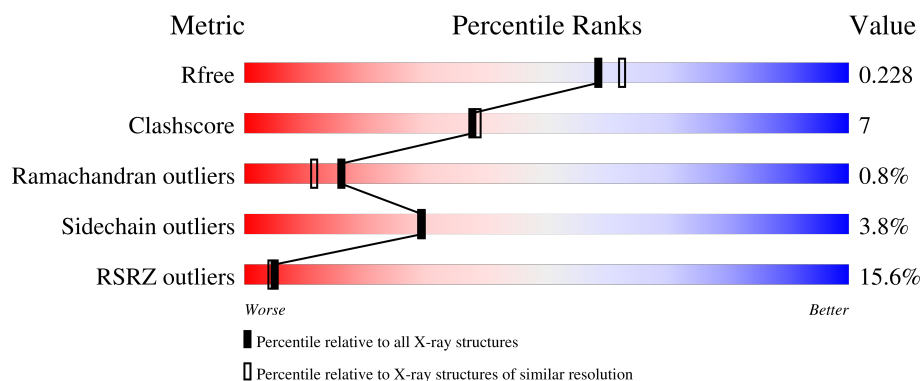
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	651	14% 80% 11% 8%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5085 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 General vesicular transport factor p115.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	S	0	0	0
			4690	2975	798	890	27			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	390	Total	O	0	0
			390	390		

3 Residue-property plots [i](#)

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 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: General vesicular transport factor p115



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	155.63Å 174.72Å 87.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.41 – 2.00 48.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	88.0 (48.41-2.00) 88.2 (48.41-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.178 , 0.204 0.204 , 0.228	Depositor DCC
R_{free} test set	7124 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5085	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.74	1/4764 (0.0%)	0.81	1/6456 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	508	ALA	CA-CB	6.75	1.63	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	518	ASN	N-CA-C	5.24	116.80	111.14

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	4690	0	4775	66	0
2	A	5	0	0	0	0
3	A	390	0	0	10	1
All	All	5085	0	4775	66	1

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 7.

All (66) 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:57:VAL:HG23	3:A:1005:HOH:O	1.43	1.16
1:A:73:ARG:HA	1:A:79:ILE:HD11	1.30	1.13
1:A:368:LEU:O	1:A:385:VAL:HG11	1.80	0.81
1:A:22:THR:OG1	3:A:1013:HOH:O	1.99	0.80
1:A:572:ILE:HD13	1:A:573:GLY:O	1.88	0.73
1:A:46:LEU:HD11	3:A:1015:HOH:O	1.90	0.72
1:A:54:ARG:O	3:A:1011:HOH:O	2.10	0.69
1:A:572:ILE:HD11	1:A:577:PHE:N	2.11	0.66
1:A:560:MET:HE3	1:A:563:LYS:HZ3	1.60	0.65
1:A:23:ILE:HG21	1:A:56:GLU:OE2	1.95	0.65
1:A:50:SER:HB3	1:A:85:THR:HG23	1.78	0.65
1:A:212:ARG:CZ	1:A:216:ILE:HD11	2.28	0.63
1:A:19:GLU:HA	1:A:22:THR:HG23	1.82	0.61
1:A:435:LEU:HD12	1:A:489:GLN:OE1	2.01	0.61
1:A:68:VAL:O	1:A:71:THR:O	2.18	0.60
1:A:139:ARG:HG2	1:A:182:ILE:HD13	1.84	0.59
1:A:319:ILE:CG1	1:A:333:THR:HG21	2.34	0.57
1:A:54:ARG:O	1:A:55:LEU:HB2	2.05	0.56
1:A:23:ILE:CG2	1:A:56:GLU:OE2	2.54	0.55
1:A:55:LEU:HD12	1:A:59:ILE:HD13	1.89	0.55
1:A:212:ARG:NH2	1:A:216:ILE:HD11	2.22	0.55
1:A:572:ILE:CD1	1:A:573:GLY:O	2.55	0.55
1:A:572:ILE:HD11	1:A:577:PHE:CA	2.37	0.54
1:A:572:ILE:HD13	1:A:572:ILE:C	2.33	0.54
1:A:73:ARG:CA	1:A:79:ILE:HD11	2.21	0.53
1:A:62:MET:HE1	1:A:116:ILE:HG22	1.91	0.53
1:A:572:ILE:CD1	1:A:577:PHE:HB2	2.39	0.53
1:A:140:TRP:CZ2	1:A:144:LYS:HD2	2.44	0.52
1:A:115:GLU:HG2	1:A:153:LEU:HD21	1.92	0.52
1:A:560:MET:HE3	1:A:563:LYS:NZ	2.24	0.52
1:A:228:VAL:HG12	3:A:750:HOH:O	2.10	0.51
1:A:133:GLU:O	1:A:139:ARG:HD3	2.10	0.50
1:A:368:LEU:HB3	1:A:385:VAL:HG13	1.93	0.50
1:A:319:ILE:HG12	1:A:333:THR:HG21	1.93	0.50
1:A:18:THR:O	1:A:22:THR:HG23	2.12	0.50
1:A:53:TYR:N	3:A:1006:HOH:O	2.44	0.49
1:A:469:ASN:HB3	3:A:899:HOH:O	2.14	0.48
1:A:49:LEU:O	1:A:53:TYR:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:THR:O	1:A:22:THR:CG2	2.63	0.47
1:A:55:LEU:O	1:A:59:ILE:HG23	2.15	0.47
1:A:321:MET:CE	1:A:356:ALA:HB2	2.45	0.47
1:A:430:PHE:CZ	1:A:463:LEU:HD11	2.50	0.47
1:A:372:MET:O	1:A:382:ARG:HD2	2.15	0.46
1:A:616:LEU:HD23	1:A:616:LEU:C	2.41	0.46
1:A:58:GLY:HA3	1:A:113:PHE:CE1	2.51	0.46
1:A:226:GLY:HA2	1:A:275:GLN:HG2	1.98	0.45
1:A:48:SER:HA	3:A:966:HOH:O	2.16	0.45
1:A:114:THR:HG22	1:A:153:LEU:HD22	1.98	0.45
1:A:55:LEU:C	1:A:55:LEU:CD1	2.89	0.45
1:A:412:ILE:O	1:A:414:ALA:N	2.50	0.45
1:A:435:LEU:HD21	1:A:604:PRO:HB3	1.99	0.45
1:A:557:GLU:O	1:A:560:MET:HE2	2.17	0.44
1:A:463:LEU:HD12	1:A:477:GLN:HG2	2.01	0.43
1:A:321:MET:HE2	1:A:356:ALA:CA	2.49	0.42
1:A:89:ILE:N	1:A:89:ILE:HD13	2.34	0.42
1:A:463:LEU:CD1	1:A:477:GLN:HG2	2.49	0.42
1:A:321:MET:HE2	1:A:356:ALA:HA	2.01	0.42
1:A:55:LEU:CD1	1:A:59:ILE:HD13	2.50	0.42
1:A:268:GLU:O	1:A:269:ASN:HB2	2.19	0.41
1:A:572:ILE:HD11	1:A:577:PHE:HB2	2.03	0.41
1:A:57:VAL:O	1:A:61:ALA:HB3	2.20	0.41
1:A:23:ILE:HG23	3:A:1015:HOH:O	2.20	0.41
1:A:181:VAL:HG23	3:A:699:HOH:O	2.20	0.41
1:A:189:LEU:C	1:A:189:LEU:HD23	2.46	0.40
1:A:23:ILE:HB	1:A:56:GLU:OE1	2.21	0.40
1:A:90:ILE:HD13	1:A:113:PHE:HB3	2.03	0.40

All (1) 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)
3:A:679:HOH:O	3:A:915:HOH:O[6_554]	2.13	0.07

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	597/651 (92%)	571 (96%)	21 (4%)	5 (1%)	16	11

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	LEU
1	A	413	ASP
1	A	269	ASN
1	A	266	GLY
1	A	268	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	532/577 (92%)	512 (96%)	20 (4%)	29	29

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	22	THR
1	A	24	GLN
1	A	35	LEU
1	A	48	SER
1	A	55	LEU

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Mol	Chain	Res	Type
1	A	56	GLU
1	A	59	ILE
1	A	89	ILE
1	A	93	ASP
1	A	139	ARG
1	A	153	LEU
1	A	312	LEU
1	A	319	ILE
1	A	323	THR
1	A	418	THR
1	A	533	GLU
1	A	566	GLN
1	A	572	ILE
1	A	575	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	70	GLN
1	A	137	HIS
1	A	476	GLN
1	A	511	HIS
1	A	518	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	A	652	-	4,4,4	0.77	0	6,6,6	1.15	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95th 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 ²)	Q<0.9
1	A	601/651 (92%)	1.00	94 (15%) 5 4	24, 33, 53, 89	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	415	THR	11.0
1	A	414	ALA	8.4
1	A	416	GLY	8.2
1	A	412	ILE	7.1
1	A	417	ASN	7.1
1	A	411	THR	7.1
1	A	56	GLU	5.8
1	A	418	THR	5.5
1	A	270	SER	5.2
1	A	110	GLY	4.9
1	A	268	GLU	4.8
1	A	632	SER	4.7
1	A	360	PRO	4.7
1	A	361	PRO	4.7
1	A	28	ASP	4.3
1	A	23	ILE	4.3
1	A	21	GLU	4.3
1	A	266	GLY	4.2
1	A	18	THR	3.9
1	A	413	ASP	3.8
1	A	410	SER	3.7
1	A	20	ALA	3.7
1	A	58	GLY	3.6
1	A	383	CYS	3.6
1	A	22	THR	3.6
1	A	529	ASN	3.6
1	A	59	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	116	ILE	3.6
1	A	435	LEU	3.5
1	A	24	GLN	3.5
1	A	52	LYS	3.5
1	A	558	THR	3.5
1	A	309	GLN	3.3
1	A	35	LEU	3.2
1	A	32	SER	3.1
1	A	271	GLY	3.1
1	A	71	THR	3.0
1	A	176	ALA	2.9
1	A	269	ASN	2.9
1	A	53	TYR	2.9
1	A	175	LEU	2.8
1	A	109	LEU	2.8
1	A	555	SER	2.8
1	A	419	VAL	2.8
1	A	359	ASN	2.8
1	A	394	TYR	2.8
1	A	49	LEU	2.8
1	A	560	MET	2.7
1	A	93	ASP	2.7
1	A	358	SER	2.6
1	A	576	ASN	2.6
1	A	530	LEU	2.6
1	A	629	TYR	2.6
1	A	553	ASP	2.6
1	A	51	LYS	2.5
1	A	459	LEU	2.5
1	A	296	PRO	2.5
1	A	54	ARG	2.5
1	A	76	SER	2.4
1	A	631	SER	2.4
1	A	357	PRO	2.4
1	A	267	ASP	2.4
1	A	386	LEU	2.4
1	A	439	CYS	2.4
1	A	445	ALA	2.4
1	A	448	LEU	2.3
1	A	169	SER	2.3
1	A	178	SER	2.3
1	A	434	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	297	PRO	2.3
1	A	324	GLY	2.3
1	A	376	ARG	2.3
1	A	531	GLY	2.3
1	A	619	GLU	2.2
1	A	295	ASN	2.2
1	A	245	SER	2.2
1	A	395	LYS	2.2
1	A	450	GLU	2.2
1	A	122	GLU	2.2
1	A	384	ALA	2.2
1	A	438	TRP	2.2
1	A	563	LYS	2.1
1	A	552	ASN	2.1
1	A	31	ALA	2.1
1	A	26	LEU	2.1
1	A	77	GLU	2.1
1	A	485	GLY	2.1
1	A	25	LYS	2.1
1	A	561	LYS	2.1
1	A	379	PHE	2.0
1	A	442	VAL	2.0
1	A	91	SER	2.0
1	A	123	ASN	2.0
1	A	162	LEU	2.0

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	A	652	5/5	0.91	0.10	67,71,73,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.