



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:08 AM GMT

PDB ID : 2W3C
Title : GLOBULAR HEAD REGION OF THE HUMAN GENERAL VESICULAR
TRANSPORT FACTOR P115
Authors : Striegl, H.; Roske, Y.; Kummel, D.; Heinemann, U.
Deposited on : 2008-11-11
Resolution : 2.22 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

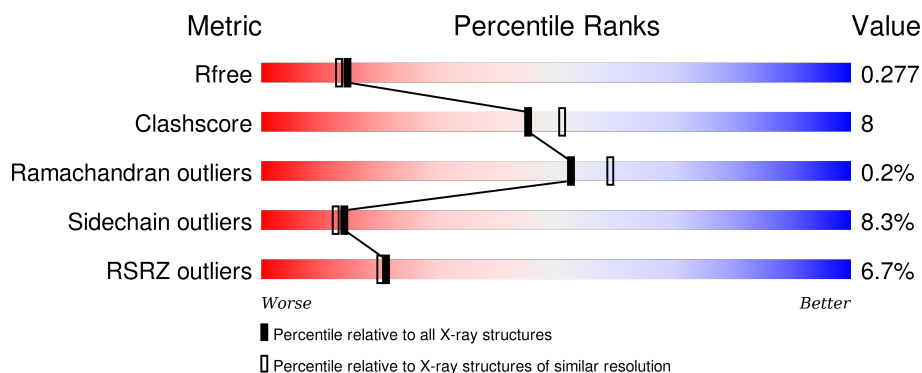
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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 ≥ 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	577	<div> <div>6%</div> <div>75%</div> <div>17%</div> <div>5%</div> </div>

2 Entry composition [i](#)

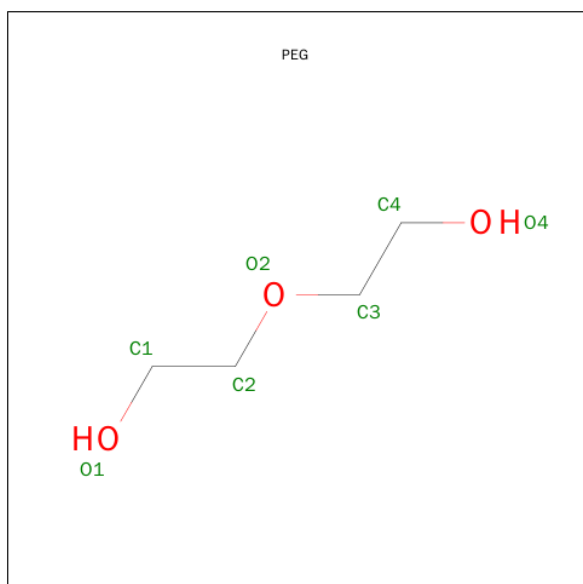
There are 3 unique types of molecules in this entry. The entry contains 4439 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
			Total	C	N	O	S			
1	A	551	4295	2737	727	805	26	0	2	0

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	7	4	3	0	0
2	A	1	7	4	3	0	0
2	A	1	7	4	3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total 123	O 123	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.55Å 68.89Å 85.75Å 90.00° 108.74° 90.00°	Depositor
Resolution (Å)	19.85 – 2.22 19.85 – 2.22	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.85-2.22) 97.9 (19.85-2.22)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.219 , 0.269 0.230 , 0.277	Depositor DCC
R_{free} test set	2352 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 62.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 47021 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4439	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.99% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PEG

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	0.57	0/4366	0.70	3/5917 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	189	LEU	CA-CB-CG	7.87	133.40	115.30
1	A	284	LEU	CA-CB-CG	6.31	129.81	115.30
1	A	288	ARG	NE-CZ-NH1	-5.30	117.65	120.30

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	4295	0	4352	68	0
2	A	21	0	30	0	0
3	A	123	0	0	4	0
All	All	4439	0	4382	68	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 8.

All (68) 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:362:ARG:HG3	1:A:362:ARG:HH11	1.16	1.09
1:A:406:THR:HG21	1:A:422:GLY:H	1.26	0.97
1:A:307:MET:HG2	1:A:312:LEU:HD23	1.48	0.92
1:A:507:ILE:H	1:A:507:ILE:HD12	1.45	0.82
1:A:406:THR:CG2	1:A:422:GLY:H	1.92	0.81
1:A:435:LEU:HG	1:A:607:MET:HE1	1.70	0.73
1:A:358:SER:HB2	1:A:360:PRO:O	1.87	0.73
1:A:212:ARG:HH11	1:A:212:ARG:HG3	1.53	0.71
1:A:362:ARG:CG	1:A:362:ARG:HH11	1.95	0.71
1:A:362:ARG:HG3	1:A:362:ARG:NH1	1.96	0.69
1:A:435:LEU:HG	1:A:607:MET:CE	2.26	0.66
1:A:212:ARG:NH1	1:A:212:ARG:HG3	2.11	0.64
1:A:624:ILE:O	1:A:628:ILE:HG12	1.98	0.62
1:A:263:PHE:HE1	1:A:284:LEU:HD13	1.64	0.61
1:A:477:GLN:O	1:A:481:ILE:HD12	2.00	0.60
1:A:263:PHE:CE1	1:A:284:LEU:HD13	2.37	0.59
1:A:115:GLU:HG2	1:A:153:LEU:HD11	1.84	0.59
1:A:406:THR:HG21	1:A:422:GLY:N	2.08	0.58
1:A:492:VAL:O	1:A:496:MET:HG3	2.06	0.55
1:A:263:PHE:CE1	1:A:284:LEU:CD1	2.89	0.55
1:A:406:THR:HG22	1:A:422:GLY:HA3	1.90	0.53
1:A:362:ARG:CG	1:A:362:ARG:NH1	2.59	0.53
1:A:77:GLU:HG2	3:A:2004:HOH:O	2.10	0.52
1:A:114:THR:HG23	1:A:149:LEU:HD23	1.92	0.52
1:A:139:ARG:O	1:A:143:VAL:HG23	2.10	0.52
1:A:87:TYR:O	1:A:91:SER:HB3	2.10	0.52
1:A:532:GLU:HA	1:A:535:GLN:HG3	1.91	0.52
1:A:406:THR:CG2	1:A:422:GLY:N	2.69	0.51
1:A:148:SER:O	1:A:152:GLN:HG3	2.11	0.51
1:A:415:THR:C	1:A:417:ASN:H	2.14	0.51
1:A:435:LEU:HD11	1:A:604:PRO:HB3	1.93	0.51
1:A:77:GLU:CG	3:A:2004:HOH:O	2.60	0.50
1:A:457:GLN:HE22	1:A:460:ARG:HH11	1.60	0.49
1:A:197:ASN:OD1	1:A:200:ILE:HG13	2.12	0.49
1:A:545:LEU:O	1:A:548:SER:HB2	2.13	0.49
1:A:139:ARG:NH1	1:A:177:ASP:OD2	2.46	0.48
1:A:549:ILE:HG12	1:A:564:LEU:HB3	1.95	0.48
1:A:183:ARG:O	1:A:187:VAL:HG23	2.12	0.48
1:A:212:ARG:HH11	1:A:212:ARG:CG	2.24	0.48
1:A:150:LEU:HB3	1:A:192:ALA:HB1	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LEU:HD21	1:A:235:LEU:HD23	1.97	0.47
1:A:403:ILE:O	1:A:406:THR:HB	2.16	0.46
1:A:507:ILE:H	1:A:507:ILE:CD1	2.21	0.46
1:A:221:GLY:O	1:A:224:ASP:HB2	2.15	0.46
1:A:423:GLN:HG3	3:A:2083:HOH:O	2.16	0.46
1:A:600:ASN:O	1:A:600:ASN:CG	2.55	0.45
1:A:375:GLU:HG2	1:A:433:ASP:OD1	2.16	0.45
1:A:488:ILE:HG13	1:A:489:GLN:N	2.31	0.45
1:A:600:ASN:C	1:A:600:ASN:ND2	2.70	0.45
1:A:186:GLY:O	1:A:189:LEU:HB3	2.15	0.45
1:A:395:LYS:HD3	1:A:395:LYS:HA	1.68	0.44
1:A:62:MET:C	1:A:64:HIS:H	2.21	0.44
1:A:527:ALA:HB2	1:A:572:ILE:HG12	1.99	0.43
1:A:364:ALA:HA	1:A:367:VAL:HG12	1.99	0.43
1:A:352:ALA:HB2	1:A:392:PHE:CE2	2.54	0.43
1:A:292:SER:OG	1:A:294:THR:HB	2.18	0.43
1:A:210:PHE:HB2	3:A:2035:HOH:O	2.18	0.43
1:A:260:LYS:HD2	1:A:311:GLY:O	2.18	0.43
1:A:127:LEU:O	1:A:131:LEU:HG	2.19	0.42
1:A:490:THR:O	1:A:494:LEU:HG	2.19	0.42
1:A:169:SER:O	1:A:172:MET:HB2	2.19	0.42
1:A:75:ASP:O	1:A:78:ILE:HG22	2.20	0.41
1:A:386:LEU:O	1:A:390:GLN:HG3	2.20	0.41
1:A:507:ILE:N	1:A:507:ILE:HD12	2.25	0.41
1:A:602:PRO:HD2	1:A:606:TYR:CG	2.55	0.41
1:A:548:SER:O	1:A:552:ASN:HB2	2.20	0.40
1:A:62:MET:C	1:A:64:HIS:N	2.74	0.40
1:A:396:ASN:O	1:A:400:GLN:HG3	2.22	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	545/577 (94%)	516 (95%)	28 (5%)	1 (0%)	52 59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	VAL

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	483/512 (94%)	443 (92%)	40 (8%)	14 12

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	VAL
1	A	72	ASP
1	A	111	SER
1	A	148	SER
1	A	150	LEU
1	A	164	SER
1	A	179	ARG
1	A	183	ARG
1	A	189	LEU
1	A	195	ARG
1	A	212	ARG
1	A	229	VAL
1	A	237	GLN
1	A	264	GLU
1	A	265	VAL
1	A	272	TRP
1	A	284	LEU
1	A	294	THR
1	A	325	VAL
1	A	362	ARG

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Mol	Chain	Res	Type
1	A	369	LEU
1	A	386	LEU
1	A	413	ASP
1	A	423	GLN
1	A	459	LEU
1	A	484	GLN
1	A	488	ILE
1	A	490	THR
1	A	491	ARG
1	A	507	ILE
1	A	513	LEU
1	A	529	ASN
1	A	533	GLU
1	A	564	LEU
1	A	567	LEU
1	A	569	GLU
1	A	580	LYS
1	A	591	SER
1	A	600	ASN
1	A	626	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	159	GLN
1	A	423	GLN
1	A	457	GLN
1	A	529	ASN
1	A	576	ASN
1	A	600	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 ligands are 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	PEG	A	1630	-	6,6,6	0.55	0	5,5,5	0.23	0
2	PEG	A	1631	-	6,6,6	0.51	0	5,5,5	0.20	0
2	PEG	A	1632	-	6,6,6	0.47	0	5,5,5	0.35	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PEG	A	1630	-	-	0/4/4/4	0/0/0/0
2	PEG	A	1631	-	-	0/4/4/4	0/0/0/0
2	PEG	A	1632	-	-	0/4/4/4	0/0/0/0

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 ⓘ

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	551/577 (95%)	0.38	37 (6%)	21 20	18, 41, 69, 84	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	LEU	6.7
1	A	146	LEU	4.6
1	A	545	LEU	4.4
1	A	471	PRO	3.9
1	A	271	GLY	3.8
1	A	57	VAL	3.6
1	A	469	ASN	3.6
1	A	467	ILE	3.5
1	A	337	VAL	3.5
1	A	340	VAL	3.3
1	A	142	GLY	3.3
1	A	388	CYS	3.3
1	A	548	SER	3.0
1	A	555	SER	3.0
1	A	470	PRO	2.9
1	A	149	LEU	2.6
1	A	361	PRO	2.6
1	A	83	LEU	2.6
1	A	92	ASN	2.6
1	A	544	LEU	2.4
1	A	134	PHE	2.4
1	A	600	ASN	2.4
1	A	250	PHE	2.4
1	A	468	GLY	2.3
1	A	549	ILE	2.3
1	A	171	LEU	2.3
1	A	492	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	341	ILE	2.2
1	A	357	PRO	2.2
1	A	262	TRP	2.2
1	A	187	VAL	2.2
1	A	531	GLY	2.1
1	A	629	TYR	2.1
1	A	391	CYS	2.1
1	A	376	ARG	2.1
1	A	203	ILE	2.0
1	A	323	THR	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	PEG	A	1630	7/7	0.87	0.13	-0.37	55,66,80,82	0
2	PEG	A	1631	7/7	0.95	0.21	-	32,45,55,60	0
2	PEG	A	1632	7/7	0.87	0.29	-	61,67,92,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.