



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:22 AM GMT

PDB ID : 3EDV
Title : Crystal Structure of Repeats 14-16 of Beta2-Spectrin
Authors : Michaely, P.; Tomchick, D.R.
Deposited on : 2008-09-03
Resolution : 1.95 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

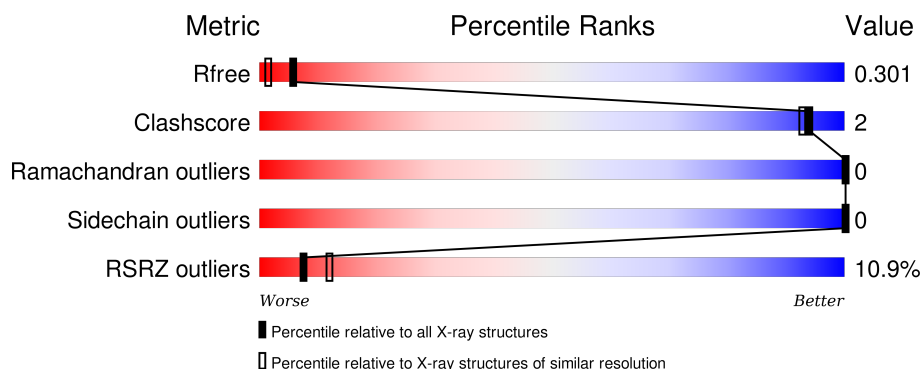
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	323	<div> <div>11%</div> <div>94%</div> <div>5%</div> </div>
1	B	323	<div> <div>11%</div> <div>97%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6186 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 Spectrin beta chain, brain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	14	0
			2753	1717	496	530	10			
1	B	321	Total	C	N	O	S	0	11	0
			2721	1692	494	524	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP Q01082
A	-3	SER	-	EXPRESSION TAG	UNP Q01082
A	-2	HIS	-	EXPRESSION TAG	UNP Q01082
A	-1	MET	-	EXPRESSION TAG	UNP Q01082
B	-4	GLY	-	EXPRESSION TAG	UNP Q01082
B	-3	SER	-	EXPRESSION TAG	UNP Q01082
B	-2	HIS	-	EXPRESSION TAG	UNP Q01082
B	-1	MET	-	EXPRESSION TAG	UNP Q01082

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	7	Total	Mg	0	0
			7	7		
2	A	6	Total	Mg	0	0
			6	6		

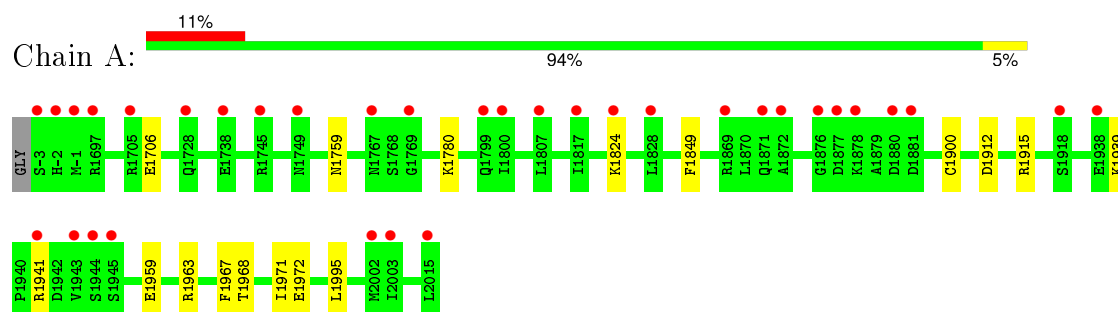
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	358	Total	O	0	0
			358	358		
3	B	341	Total	O	0	0
			341	341		

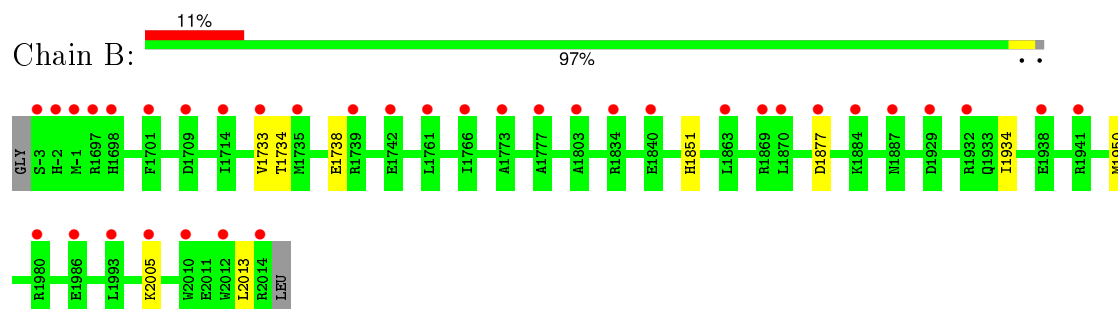
3 Residue-property plots

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.

- Molecule 1: Spectrin beta chain, brain 1



- Molecule 1: Spectrin beta chain, brain 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.34Å 56.41Å 261.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.63 – 1.95 28.04 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.2 (28.63-1.95) 97.2 (28.04-1.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.192 , 0.250 0.257 , 0.301	Depositor DCC
R_{free} test set	2788 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	29.8	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 54964 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6186	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.82% 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: MG

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	0.64	1/2828 (0.0%)	0.59	0/3802
1	B	0.58	0/2799	0.60	0/3764
All	All	0.61	1/5627 (0.0%)	0.60	0/7566

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1900	CYS	CB-SG	-5.93	1.72	1.81

There are no bond angle outliers.

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	2753	0	2715	14	0
1	B	2721	0	2671	7	0
2	A	6	0	0	0	0
2	B	7	0	0	0	0
3	A	358	0	0	0	0
3	B	341	0	0	1	0
All	All	6186	0	5386	20	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 2.

All (20) 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:1971[B]:ILE:HD11	1:A:1995:LEU:HD23	1.39	1.03
1:A:1971[B]:ILE:HD11	1:A:1995:LEU:CD2	2.12	0.79
1:A:1971[B]:ILE:CD1	1:A:1995:LEU:HD23	2.12	0.77
1:B:1950[B]:MET:HE2	1:B:2013:LEU:HB3	1.88	0.55
1:A:1967:PHE:O	1:A:1971[B]:ILE:HG12	2.08	0.53
1:A:1971[B]:ILE:CD1	1:A:1995:LEU:CD2	2.83	0.51
1:A:1759:ASN:OD1	1:A:1780[B]:LYS:NZ	2.33	0.50
1:A:1939[B]:LYS:HE2	1:A:1941:ARG:NH1	2.27	0.50
1:B:1877[B]:ASP:HB2	3:B:366:HOH:O	2.11	0.49
1:A:1939[B]:LYS:HB3	1:A:1939[B]:LYS:NZ	2.30	0.46
1:B:1950[B]:MET:CE	1:B:2013:LEU:HB3	2.45	0.46
1:A:1824[B]:LYS:HE2	1:A:1849:PHE:CE1	2.50	0.45
1:A:1824[B]:LYS:HD2	1:A:1824[B]:LYS:HA	1.69	0.44
1:A:1706:GLU:HG3	1:B:1851:HIS:NE2	2.32	0.44
1:A:1968:THR:O	1:A:1972[B]:GLU:HG2	2.19	0.43
1:B:1733:VAL:HG13	1:B:1734:THR:N	2.34	0.42
1:A:1959[B]:GLU:HG3	1:A:1963:ARG:HH12	1.85	0.41
1:A:1912:ASP:OD1	1:A:1915:ARG:NH1	2.49	0.41
1:B:1934:ILE:HG21	1:B:2005[B]:LYS:HG3	2.02	0.40
1:B:1738:GLU:OE1	1:B:1738:GLU:HA	2.21	0.40

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	334/323 (103%)	333 (100%)	1 (0%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	330/323 (102%)	329 (100%)	1 (0%)	0	100	100
All	All	664/646 (103%)	662 (100%)	2 (0%)	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	297/283 (105%)	297 (100%)	0	100	100
1	B	293/283 (104%)	293 (100%)	0	100	100
All	All	590/566 (104%)	590 (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. There are no such sidechains identified.

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](#)

Of 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

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	322/323 (99%)	0.93	34 (10%) 8 13	7, 23, 32, 36	5 (1%)
1	B	321/323 (99%)	0.97	36 (11%) 7 11	11, 27, 37, 44	6 (1%)
All	All	643/646 (99%)	0.95	70 (10%) 7 12	7, 26, 35, 44	11 (1%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-3	SER	7.9
1	A	-2	HIS	6.7
1	B	-3	SER	6.2
1	B	1701	PHE	5.8
1	B	1735	MET	5.7
1	A	1877	ASP	5.6
1	B	1877[A]	ASP	4.8
1	B	1834	ARG	4.6
1	A	1876	GLY	4.2
1	B	-2	HIS	4.1
1	B	1733	VAL	4.1
1	B	1993	LEU	4.0
1	A	1943	VAL	3.6
1	A	1807	LEU	3.5
1	A	1800	ILE	3.4
1	B	1980[A]	ARG	3.3
1	A	1728	GLN	3.3
1	B	2014	ARG	3.3
1	B	1840	GLU	3.2
1	A	1944	SER	3.1
1	B	2010	TRP	3.1
1	A	1880	ASP	3.1
1	A	2015	LEU	3.0
1	B	1869[A]	ARG	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	1918[A]	SER	3.0
1	B	2005[A]	LYS	3.0
1	A	1749	ASN	2.9
1	A	2003[A]	ILE	2.9
1	B	1697	ARG	2.8
1	A	1705	ARG	2.8
1	A	1878	LYS	2.8
1	A	1828	LEU	2.8
1	B	1941	ARG	2.7
1	B	1739	ARG	2.7
1	B	1709[A]	ASP	2.7
1	A	1799	GLN	2.6
1	A	1745	ARG	2.6
1	B	1884	LYS	2.5
1	B	1698	HIS	2.5
1	A	1941	ARG	2.5
1	A	1824[A]	LYS	2.4
1	B	1887	ASN	2.4
1	B	1863	LEU	2.4
1	A	1767	ASN	2.4
1	B	1803	ALA	2.4
1	B	1742	GLU	2.3
1	A	1738	GLU	2.3
1	B	1986	GLU	2.3
1	A	-1	MET	2.3
1	B	1932	ARG	2.3
1	A	1769	GLY	2.3
1	A	1872	ALA	2.3
1	B	1766	ILE	2.2
1	A	1881	ASP	2.2
1	B	1929	ASP	2.2
1	A	1938	GLU	2.2
1	B	1870	LEU	2.2
1	A	2002	MET	2.2
1	A	1871	GLN	2.1
1	A	1945	SER	2.1
1	B	1938	GLU	2.1
1	B	1773	ALA	2.1
1	A	1697	ARG	2.1
1	B	-1	MET	2.1
1	B	2012	TRP	2.1
1	B	1761	LEU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	1817	ILE	2.0
1	A	1869[A]	ARG	2.0
1	B	1777	ALA	2.0
1	B	1714	ILE	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(Å ²)	Q<0.9
2	MG	A	1007	1/1	0.82	0.18	-0.42	30,30,30,30	0
2	MG	B	1002	1/1	0.88	0.23	-	73,73,73,73	0
2	MG	B	1009	1/1	0.79	0.20	-	30,30,30,30	0
2	MG	B	1012	1/1	0.84	0.17	-	30,30,30,30	0
2	MG	B	1010	1/1	0.84	0.20	-	30,30,30,30	0
2	MG	A	1004	1/1	0.76	0.17	-	30,30,30,30	0
2	MG	B	1003	1/1	0.33	0.20	-	30,30,30,30	0
2	MG	A	1006	1/1	0.94	0.26	-	30,30,30,30	0
2	MG	A	1011	1/1	0.75	0.53	-	30,30,30,30	0
2	MG	A	1005	1/1	0.43	0.27	-	30,30,30,30	0
2	MG	B	1000	1/1	0.91	0.18	-	62,62,62,62	0
2	MG	B	1001	1/1	0.95	0.10	-	56,56,56,56	0
2	MG	A	1008	1/1	0.47	0.19	-	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.