



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 09:38 pm GMT

PDB ID : 2FFB
Title : The crystal structure of the HlyB-NBD E631Q mutant in complex with ADP
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Deposited on : 2005-12-19
Resolution : 1.90 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

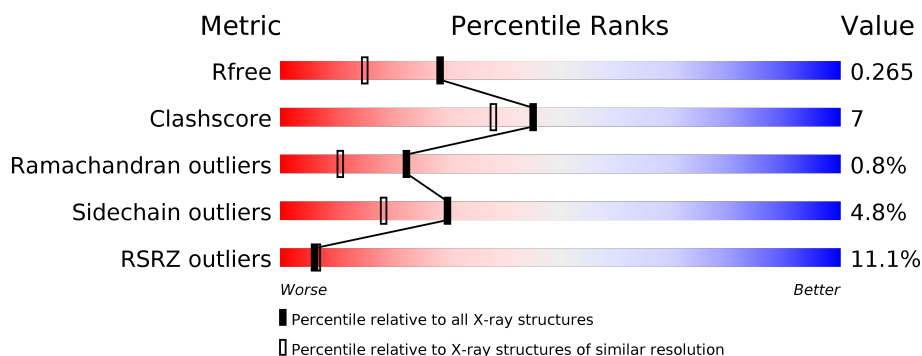
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

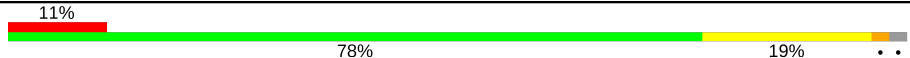
The reported resolution of this entry is 1.90 Å.

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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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	247	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2067 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.

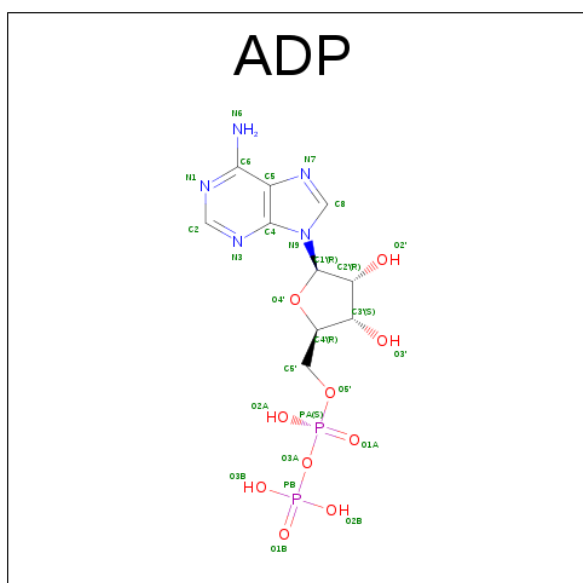
- Molecule 1 is a protein called Alpha-hemolysin translocation ATP-binding protein hlyB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1912	1206	348	353	5			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	461	HIS	-	EXPRESSION TAG	UNP P08716
A	462	HIS	-	EXPRESSION TAG	UNP P08716
A	463	HIS	-	EXPRESSION TAG	UNP P08716
A	464	HIS	-	EXPRESSION TAG	UNP P08716
A	465	HIS	-	EXPRESSION TAG	UNP P08716
A	466	HIS	-	EXPRESSION TAG	UNP P08716
A	631	GLN	GLU	ENGINEERED	UNP P08716

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	128	Total	O	0	0
			128	128		

i

- Molecule 1: Alpha-hemolysin translocation ATP-binding protein hlyB



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.57Å 34.75Å 37.52Å 90.00° 97.97° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 19.75 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.1 (20.00-1.90) 94.2 (19.75-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.202 , 0.250 0.214 , 0.265	Depositor DCC
R_{free} test set	876 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2067	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.28% 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: ADP

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	0.58	0/1941	0.80	3/2619 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	480	ASP	CB-CG-OD2	5.59	123.33	118.30
1	A	637	ASP	CB-CG-OD2	5.49	123.25	118.30
1	A	707	ASP	CB-CG-OD2	5.07	122.86	118.30

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	1912	0	1963	29	0
2	A	27	0	12	0	0
3	A	128	0	0	8	0
All	All	2067	0	1975	29	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 7.

All (29) 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:631:GLN:HE21	1:A:662:HIS:HD2	1.23	0.82
1:A:496:GLU:OE1	1:A:672:ARG:HD3	1.80	0.80
1:A:623:ASN:HD22	1:A:655:ARG:HH12	1.27	0.79
1:A:693:GLU:HB2	1:A:696:SER:HB3	1.66	0.76
1:A:631:GLN:HE21	1:A:662:HIS:CD2	2.03	0.76
1:A:466:HIS:CD2	1:A:491:SER:HB3	2.25	0.70
1:A:650:LYS:HD2	3:A:827:HOH:O	1.93	0.66
1:A:668:LYS:HD3	3:A:763:HOH:O	1.98	0.64
1:A:647:ASN:CG	3:A:827:HOH:O	2.37	0.62
1:A:529:ASP:OD2	1:A:625:LYS:NZ	2.32	0.58
1:A:623:ASN:ND2	1:A:655:ARG:HH12	1.99	0.58
1:A:531:HIS:ND1	3:A:746:HOH:O	2.34	0.53
1:A:636:LEU:HB2	1:A:641:GLU:HG2	1.91	0.52
1:A:572:VAL:O	1:A:576:ILE:HG12	2.11	0.50
1:A:647:ASN:ND2	3:A:827:HOH:O	2.43	0.50
1:A:631:GLN:NE2	1:A:662:HIS:HD2	2.02	0.50
1:A:700:TYR:CE2	1:A:704:LEU:HD11	2.50	0.47
1:A:601:GLU:O	1:A:602:GLN:HB3	2.16	0.46
1:A:501:VAL:HG11	1:A:664:LEU:HD22	1.96	0.46
1:A:697:LEU:HG	1:A:701:LEU:CD2	2.47	0.44
1:A:567:ASN:O	1:A:570:MET:HG2	2.17	0.44
1:A:470:PHE:CE2	1:A:511:LEU:HD12	2.53	0.44
1:A:475:PHE:CE1	1:A:519:TYR:HB3	2.54	0.43
1:A:631:GLN:NE2	1:A:662:HIS:CD2	2.79	0.42
1:A:693:GLU:CG	3:A:717:HOH:O	2.68	0.42
1:A:484:ILE:HG22	1:A:485:LEU:HD13	2.02	0.41
1:A:635:ALA:HA	3:A:798:HOH:O	2.21	0.41
1:A:577:TYR:CE1	3:A:772:HOH:O	2.71	0.41
1:A:526:VAL:C	1:A:527:LEU:HD12	2.41	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	241/247 (98%)	233 (97%)	6 (2%)	2 (1%)	22	11

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	566	ALA
1	A	604	ALA

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	209/213 (98%)	199 (95%)	10 (5%)	30	18

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	485	LEU
1	A	517	ARG
1	A	544	GLN
1	A	567	ASN
1	A	587	PHE
1	A	591	LEU
1	A	646	ARG
1	A	647	ASN
1	A	663	ARG
1	A	701	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	466	HIS
1	A	531	HIS

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Mol	Chain	Res	Type
1	A	552	ASN
1	A	567	ASN
1	A	610	GLN
1	A	612	GLN
1	A	623	ASN
1	A	642	HIS
1	A	647	ASN
1	A	662	HIS

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](#)

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	ADP	A	708	-	25,29,29	1.24	3 (12%)	24,45,45	2.51	3 (12%)

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	ADP	A	708	-	-	0/12/32/32	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	708	ADP	O4'-C4'	-2.10	1.40	1.45
2	A	708	ADP	C2-N1	2.88	1.39	1.33
2	A	708	ADP	C2-N3	3.37	1.37	1.32

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	708	ADP	N3-C2-N1	-11.52	118.82	128.86
2	A	708	ADP	C1'-N9-C4	-2.28	122.70	126.64
2	A	708	ADP	O3B-PB-O2B	2.33	117.03	107.61

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	243/247 (98%)	0.52	27 (11%) 6 6	21, 38, 66, 107	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	604	ALA	8.4
1	A	601	GLU	7.6
1	A	569	GLY	5.3
1	A	602	GLN	4.6
1	A	600	GLY	4.3
1	A	480	ASP	4.1
1	A	614	ILE	4.0
1	A	707	ASP	4.0
1	A	559	ILE	3.8
1	A	563	ILE	3.7
1	A	633	THR	3.2
1	A	603	GLY	3.2
1	A	515	ILE	3.0
1	A	511	LEU	2.9
1	A	588	ILE	2.9
1	A	566	ALA	2.7
1	A	616	ILE	2.7
1	A	526	VAL	2.7
1	A	617	ALA	2.6
1	A	635	ALA	2.5
1	A	634	SER	2.5
1	A	662	HIS	2.3
1	A	694	PRO	2.2
1	A	565	LEU	2.1
1	A	539	ASN	2.1
1	A	706	SER	2.0
1	A	618	ARG	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	ADP	A	708	27/27	0.98	0.07	-1.21	33,39,43,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.