



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 03:59 am BST

PDB ID : 2UZ2  
Title : Crystal structure of Xenavidin  
Authors : Helppolainen, S.H.; Maatta, J.A.E.; Airene, T.T.; Johnson, M.S.; Kulomaa, M.S.; Nordlund, H.R.  
Deposited on : 2007-04-24  
Resolution : 1.70 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

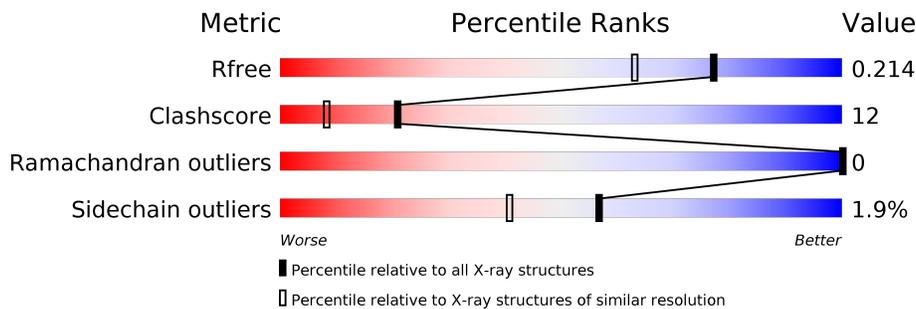
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)

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  $\geq 3$ , 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	130	 67%                      22%                      12%
1	D	130	 76%                      15%                      8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	1123	-	-	X	-

## 2 Entry composition [i](#)

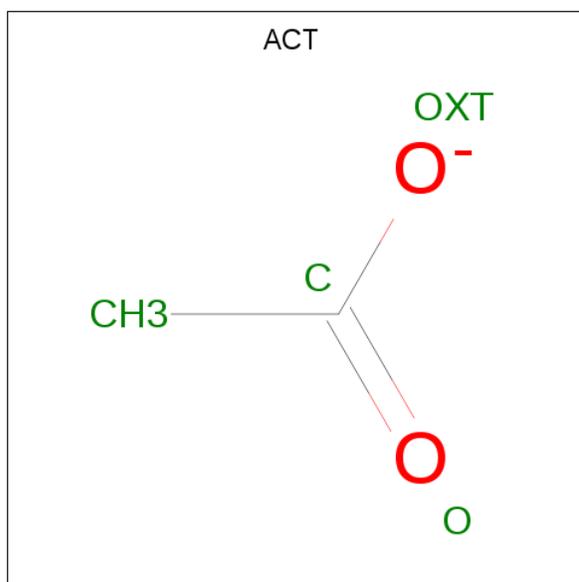
There are 4 unique types of molecules in this entry. The entry contains 2225 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 XENAVIDIN.

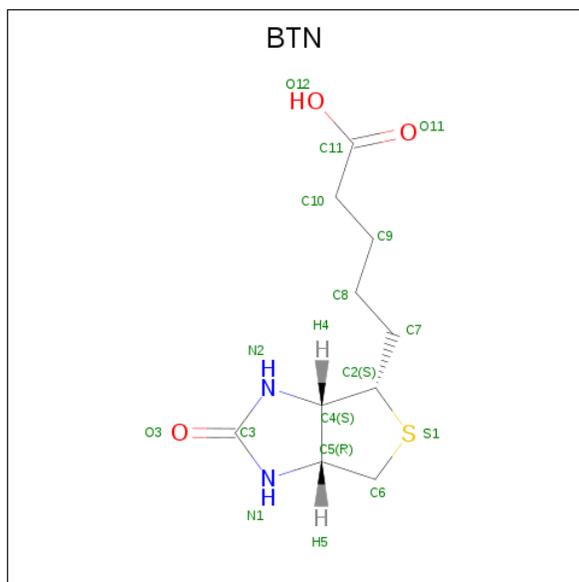
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	115	Total 972	C 621	N 161	O 186	S 4	0	9	0
1	D	119	Total 1025	C 651	N 171	O 199	S 4	0	15	0

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).



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

- Molecule 3 is BIOTIN (three-letter code: BTN) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>2</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	D	1	16	10	2	3	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	96	Total	O	0	0
			96	96		
4	D	112	Total	O	0	0
			112	112		



## 4 Data and refinement statistics i

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.82Å 110.82Å 142.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 1.70 24.93 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (25.00-1.70) 99.9 (24.93-1.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.86 (at 1.70Å)	Xtrriage
Refinement program	REFMAC 5.3.0021	Depositor
R, $R_{free}$	0.181 , 0.216 0.182 , 0.214	Depositor DCC
$R_{free}$ test set	1852 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.1	Xtrriage
Anisotropy	0.635	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 33.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.027 for $-2/3^*h-1/3^*k+2/3^*l,-1/3^*h-2/3^*k-2/3^*l,2/3^*h-2/3^*k+1/3^*l$ 0.019 for $-h,1/3^*h-1/3^*k+2/3^*l,2/3^*h+4/3^*k+1/3^*l$ 0.017 for $-1/3^*h+1/3^*k-2/3^*l,-k,-4/3^*h-2/3^*k+1/3^*l$	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2225	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	5.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.66	0/1017	0.71	0/1376
1	D	0.77	1/1086 (0.1%)	0.80	1/1472 (0.1%)
All	All	0.72	1/2103 (0.0%)	0.76	1/2848 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	54	TYR	CD2-CE2	-5.53	1.31	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	54	TYR	CB-CG-CD2	-5.80	117.52	121.00

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	972	0	963	29	0
1	D	1025	0	1027	24	0
2	A	4	0	3	2	0
3	D	16	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	96	0	0	6	0
4	D	112	0	0	5	0
All	All	2225	0	2008	47	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 12.

All (47) 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:23:GLU:OE2	4:A:2012:HOH:O	2.01	0.77
2:A:1123:ACT:H1	4:A:2089:HOH:O	1.89	0.72
1:A:105[B]:GLU:OE2	1:D:91[B]:ILE:HD11	1.93	0.69
1:A:4:GLN:NE2	1:A:88:GLY:O	2.25	0.69
1:D:91[A]:ILE:CG2	1:D:118[A]:THR:CG2	2.73	0.67
2:A:1123:ACT:H3	4:A:2023:HOH:O	1.96	0.65
1:A:6:CYS:CB	1:A:61:PRO:HD2	2.28	0.63
1:D:91[A]:ILE:CG2	1:D:118[A]:THR:HG23	2.29	0.63
1:D:40[B]:SER:OG	1:D:42:THR:HG22	2.01	0.60
1:D:91[B]:ILE:HG22	1:D:120:THR:HG22	1.83	0.60
1:A:91[B]:ILE:HD11	1:D:105:GLU:OE2	2.02	0.59
1:D:19[A]:ASN:ND2	4:D:2011:HOH:O	2.30	0.59
1:A:36:PHE:HE1	1:A:46:ILE:HG23	1.68	0.58
1:A:91[B]:ILE:HG22	1:A:120:THR:HG22	1.84	0.58
1:D:91[A]:ILE:HG23	1:D:118[A]:THR:CG2	2.34	0.57
1:A:109:TRP:CZ2	1:A:110[B]:THR:HG22	2.39	0.57
1:A:32[A]:THR:HG22	4:A:2014:HOH:O	2.06	0.56
1:D:59:GLU:HG2	1:D:60:LYS:HG3	1.88	0.55
1:A:32[B]:THR:HG22	4:D:2028:HOH:O	2.08	0.53
1:A:6:CYS:HB2	1:A:61:PRO:HD2	1.90	0.53
1:A:32[B]:THR:CG2	4:D:2028:HOH:O	2.57	0.52
1:D:107:ASP:O	1:D:110[B]:THR:HG22	2.09	0.52
1:D:32[A]:THR:HG22	4:D:2053:HOH:O	2.11	0.51
1:A:52:THR:HG22	1:D:54:TYR:CE2	2.45	0.51
1:D:118[B]:THR:HG21	4:D:2111:HOH:O	2.10	0.50
1:A:4:GLN:HE22	1:A:89:GLU:HA	1.76	0.49
1:A:105[B]:GLU:OE2	1:D:91[B]:ILE:CD1	2.58	0.49
1:D:85:ASN:HD22	1:D:91[A]:ILE:HD11	1.76	0.49
1:A:109:TRP:CE2	1:A:110[B]:THR:CG2	2.95	0.49
1:A:107:ASP:O	1:A:110[A]:THR:HG22	2.12	0.48
1:D:104:LYS:HG3	1:D:106:GLN:HG2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2028:HOH:O	1:D:52[B]:THR:CG2	2.62	0.47
1:A:105[B]:GLU:CD	1:D:91[B]:ILE:HD11	2.36	0.45
1:A:6:CYS:HB2	1:A:61:PRO:CD	2.46	0.45
1:D:91[A]:ILE:HG22	1:D:118[A]:THR:HG23	1.99	0.45
1:A:109:TRP:CE2	1:A:110[B]:THR:HG22	2.53	0.43
4:A:2028:HOH:O	1:D:52[B]:THR:HG22	2.18	0.42
1:A:54:TYR:OH	1:D:52[B]:THR:HG23	2.19	0.42
1:A:54:TYR:CG	1:D:68:HIS:HB2	2.54	0.42
1:A:56:LYS:HE2	1:D:73:ASP:HA	2.01	0.41
1:A:6:CYS:CB	1:A:82:CYS:SG	3.08	0.41
1:A:95[B]:MET:SD	1:A:116:ALA:HB2	2.61	0.41
1:A:16:LEU:HD22	1:A:39:VAL:HG11	2.03	0.40
1:A:109:TRP:CH2	1:A:110[B]:THR:HG22	2.56	0.40
1:A:5:LYS:HD2	1:A:5:LYS:HA	1.82	0.40
1:D:109:TRP:CE2	1:D:110[A]:THR:HG22	2.57	0.40
1:A:109:TRP:CZ2	1:A:110[B]:THR:CG2	3.05	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	120/130 (92%)	117 (98%)	3 (2%)	0	100	100
1	D	132/130 (102%)	130 (98%)	2 (2%)	0	100	100
All	All	252/260 (97%)	247 (98%)	5 (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	112/117 (96%)	108 (96%)	4 (4%)	35	16
1	D	122/117 (104%)	122 (100%)	0	100	100
All	All	234/234 (100%)	230 (98%)	4 (2%)	57	46

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	SER
1	A	68	HIS
1	A	72	SER
1	A	87	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	D	11	GLN

### 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

2 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	ACT	A	1123	-	1,3,3	0.82	0	0,3,3	0.00	-
3	BTN	D	1124	-	14,17,17	3.44	2 (14%)	19,23,23	2.51	6 (31%)

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
3	BTN	D	1124	-	-	1/5/28/28	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1124	BTN	C2-S1	-9.46	1.67	1.82
3	D	1124	BTN	O3-C3	8.47	1.41	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1124	BTN	C5-C6-S1	5.90	111.36	106.31
3	D	1124	BTN	C2-C4-N2	5.19	117.78	113.13
3	D	1124	BTN	C4-C2-S1	4.88	109.85	105.20
3	D	1124	BTN	C6-C5-C4	-3.08	105.99	108.66
3	D	1124	BTN	C2-C4-C5	2.22	111.52	108.94
3	D	1124	BTN	C4-N2-C3	-2.20	110.57	112.62

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1124	BTN	C11-C10-C9-C8

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1123	ACT	2	0

## 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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.