



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 14, 2017 – 11:52 pm GMT

PDB ID : 3V4R  
Title : Crystal structure of a UvrB dimer-DNA complex  
Authors : Webster, M.P.J.; Jukes, R.; Barrett, T.  
Deposited on : 2011-12-15  
Resolution : 3.25 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<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

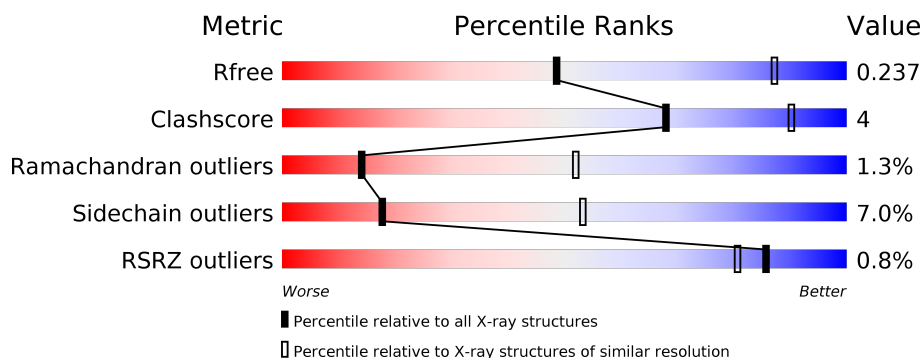
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.25 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	1852 (3.32-3.20)
Clashscore	112137	2036 (3.32-3.20)
Ramachandran outliers	110173	2000 (3.32-3.20)
Sidechain outliers	110143	1998 (3.32-3.20)
RSRZ outliers	101464	1861 (3.32-3.20)

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  $\geq 3$ , 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	667	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 75%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>75%</span> <span>12%</span> <span>•</span> <span>12%</span> </div> </div>
1	B	667	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 73%, yellow 12%, orange 1%, grey 13%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>73%</span> <span>12%</span> <span>•</span> <span>13%</span> </div> </div>
2	C	8	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 13%, yellow 25%, orange 50%, grey 13%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>13%</span> <span>25%</span> <span>50%</span> <span>13%</span> </div> </div>
2	D	8	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 13%, yellow 63%, orange 13%, grey 13%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>13%</span> <span>63%</span> <span>13%</span> <span>13%</span> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8815 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 UvrABC system protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	586	Total	C	N	O	S	0	0	0
			4321	2738	754	815	14			
1	B	579	Total	C	N	O	S	0	0	0
			4205	2663	739	790	13			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP P37954
A	-4	HIS	-	EXPRESSION TAG	UNP P37954
A	-3	HIS	-	EXPRESSION TAG	UNP P37954
A	-2	HIS	-	EXPRESSION TAG	UNP P37954
A	-1	HIS	-	EXPRESSION TAG	UNP P37954
A	0	HIS	-	EXPRESSION TAG	UNP P37954
B	-5	HIS	-	EXPRESSION TAG	UNP P37954
B	-4	HIS	-	EXPRESSION TAG	UNP P37954
B	-3	HIS	-	EXPRESSION TAG	UNP P37954
B	-2	HIS	-	EXPRESSION TAG	UNP P37954
B	-1	HIS	-	EXPRESSION TAG	UNP P37954
B	0	HIS	-	EXPRESSION TAG	UNP P37954

- Molecule 2 is a DNA chain called DNA: 5'-TACTGTTT-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	7	Total	C	N	O	P	0	0	0
			131	64	20	41	6			
2	D	7	Total	C	N	O	P	0	0	0
			123	59	19	39	6			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).

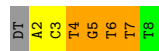


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	0	0	
			27	10	5	10	2			
3	B	1	Total	O	P				0	0
			8	6	2					




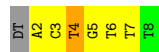
- Molecule 2: DNA: 5 -TACTGTTT-3

Chain C: 



- Molecule 2: DNA: 5 -TACTGTTT-3

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.94Å 100.31Å 163.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.91 – 3.25 19.91 – 3.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.91-3.25) 100.0 (19.91-3.25)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.93 (at 3.22Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.180 , 0.219 0.192 , 0.237	Depositor DCC
$R_{free}$ test set	1297 reflections (5.11%)	DCC
Wilson B-factor (Å <sup>2</sup> )	75.2	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 65.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.027 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8815	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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: 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	0.52	0/4397	0.71	0/5996
1	B	0.50	0/4278	0.74	3/5843 (0.1%)
2	C	1.19	0/145	2.40	14/222 (6.3%)
2	D	1.22	0/136	2.50	12/209 (5.7%)
All	All	0.54	0/8956	0.85	29/12270 (0.2%)

There are no bond length outliers.

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	6	DT	O4'-C1'-N1	9.27	114.49	108.00
2	C	2	DA	O4'-C1'-N9	8.97	114.28	108.00
2	D	6	DT	O4'-C1'-N1	8.93	114.25	108.00
2	D	7	DT	C4-C5-C7	8.81	124.29	119.00
2	D	6	DT	P-O3'-C3'	8.37	129.75	119.70

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	4321	0	3986	38	0
1	B	4205	0	3844	35	0

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	131	0	78	3	0
2	D	123	0	70	1	0
3	A	27	0	12	0	0
3	B	8	0	0	0	0
All	All	8815	0	7990	74	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 4.

The worst 5 of 74 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)
2:C:4:DT:H2''	2:C:5:DG:H5'	1.51	0.90
2:D:4:DT:H2''	2:D:5:DG:H5''	1.63	0.79
1:A:181:TYR:CE2	1:A:246:ALA:HB2	2.17	0.78
1:A:51:ASN:O	1:A:55:GLU:HG2	1.94	0.67
1:B:32:LYS:HA	1:B:33:LYS:HB2	1.79	0.64

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	582/667 (87%)	546 (94%)	30 (5%)	6 (1%)	18	58
1	B	575/667 (86%)	543 (94%)	23 (4%)	9 (2%)	11	47
All	All	1157/1334 (87%)	1089 (94%)	53 (5%)	15 (1%)	14	52

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	209	GLU
1	A	587	THR
1	B	33	LYS
1	B	163	THR
1	B	165	MET

### 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	413/592 (70%)	386 (94%)	27 (6%)	20	56
1	B	392/592 (66%)	363 (93%)	29 (7%)	16	50
All	All	805/1184 (68%)	749 (93%)	56 (7%)	18	53

5 of 56 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	510	ASP
1	B	92	TYR
1	B	475	LEU
1	A	583	ILE
1	B	60	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	355	GLN
1	B	439	GLN
1	B	359	GLN
1	A	364	HIS
1	B	364	HIS

### 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 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	$\# Z  > 2$
3	ADP	A	701	-	25,29,29	0.60	0	24,45,45	0.93	1 (4%)
3	ADP	B	701	-	4,7,29	0.42	0	4,10,45	0.40	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
3	ADP	A	701	-	-	0/12/32/32	0/3/3/3
3	ADP	B	701	-	-	0/2/5/32	0/0/0/3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	ADP	O5'-C5'-C4'	2.54	118.02	109.00

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

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	586/667 (87%)	-0.51	6 (1%) 82 76	32, 65, 94, 128	0
1	B	579/667 (86%)	-0.47	3 (0%) 90 87	36, 69, 98, 113	0
2	C	7/8 (87%)	0.17	0 100 100	79, 89, 95, 109	0
2	D	7/8 (87%)	-0.02	0 100 100	72, 75, 85, 88	0
All	All	1179/1350 (87%)	-0.49	9 (0%) 86 80	32, 68, 97, 128	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	589	ASN	3.3
1	A	587	THR	3.0
1	B	29	GLN	2.9
1	B	239	ASP	2.7
1	A	248	HIS	2.7

### 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	ADP	A	701	27/27	0.87	0.27	1.90	100,105,109,110	0
3	ADP	B	701	8/27	0.94	0.13	-0.69	117,119,121,122	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.