



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:19 PM UTC

PDB ID : 1KT6 / pdb_00001kt6
Title : Crystal structure of bovine holo-RBP at pH 9.0
Authors : Calderone, V.; Berni, R.; Zanotti, G.
Deposited on : 2002-01-15
Resolution : 1.10 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

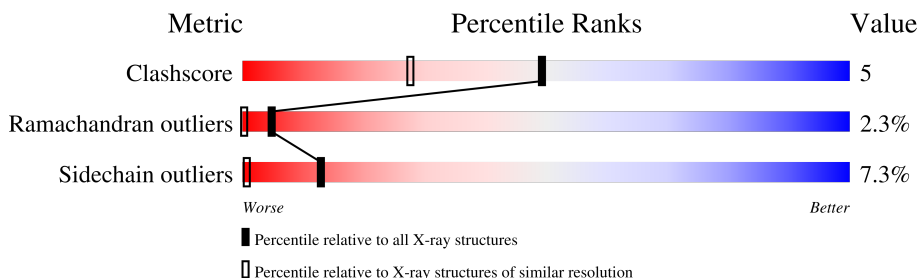
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.10 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))
Clashscore	190562	1656 (1.12-1.08)
Ramachandran outliers	187476	1616 (1.12-1.08)
Sidechain outliers	187428	1612 (1.12-1.08)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	183	 75% 11% 5% . .

2 Entry composition [i](#)

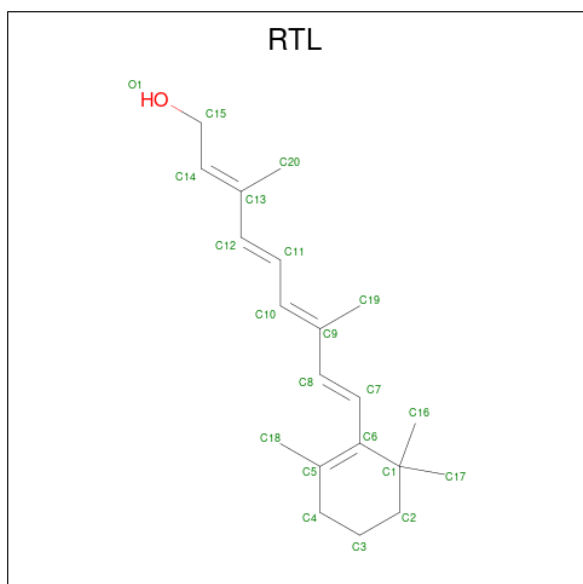
There are 3 unique types of molecules in this entry. The entry contains 1611 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 plasma retinol-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	175	1418	893	246	269	10	0	0	0

- Molecule 2 is RETINOL (CCD ID: RTL) (formula: C₂₀H₃₀O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	21	20	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	172	Total	O	0	0
			172	172		

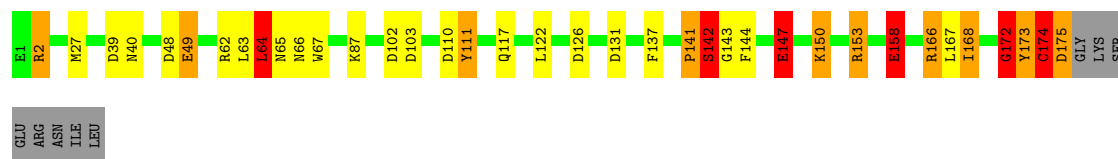
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

- Molecule 1: plasma retinol-binding protein

Chain A: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.86 Å 48.95 Å 75.65 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.42 – 1.10	Depositor
% Data completeness (in resolution range)	83.4 (22.42-1.10)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.188 , 0.223	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1611	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.11	6/1452 (0.4%)	2.15	37/1961 (1.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	143	GLY	C-N	14.54	1.54	1.33
1	A	173	TYR	N-CA	-13.93	1.28	1.46
1	A	111	TYR	N-CA	-13.48	1.28	1.46
1	A	172	GLY	C-N	-11.43	1.17	1.33
1	A	144	PHE	N-CA	9.78	1.58	1.46
1	A	110	ASP	C-N	-6.20	1.25	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	172	GLY	O-C-N	-26.70	87.99	122.70
1	A	172	GLY	CA-C-N	26.61	172.37	121.54
1	A	172	GLY	C-N-CA	26.61	172.37	121.54
1	A	110	ASP	CA-C-N	24.09	167.56	121.54
1	A	110	ASP	C-N-CA	24.09	167.56	121.54
1	A	110	ASP	O-C-N	-16.91	102.12	122.25
1	A	143	GLY	O-C-N	11.72	135.85	122.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	GLU	CB-CG-CD	9.30	128.42	112.60
1	A	142	SER	CA-C-N	9.16	140.50	121.17
1	A	142	SER	C-N-CA	9.16	140.50	121.17
1	A	143	GLY	CA-C-N	-8.39	105.52	121.54
1	A	143	GLY	C-N-CA	-8.39	105.52	121.54
1	A	144	PHE	N-CA-CB	7.90	123.83	110.49
1	A	174	CYS	CA-C-N	7.87	135.87	121.70
1	A	174	CYS	C-N-CA	7.87	135.87	121.70
1	A	49	GLU	CB-CG-CD	7.77	125.82	112.60
1	A	147	GLU	CB-CG-CD	7.55	125.44	112.60
1	A	27	MET	CG-SD-CE	7.36	117.10	100.90
1	A	173	TYR	CA-CB-CG	-6.84	101.59	113.90
1	A	153	ARG	NE-CZ-NH2	6.71	125.24	119.20
1	A	64	LEU	CA-C-N	6.36	132.78	122.67
1	A	64	LEU	C-N-CA	6.36	132.78	122.67
1	A	141	PRO	O-C-N	6.36	131.22	122.64
1	A	166	ARG	CA-CB-CG	6.16	126.41	114.10
1	A	126	ASP	CA-CB-CG	6.00	118.60	112.60
1	A	168	ILE	CA-CB-CG1	5.84	120.34	110.40
1	A	137	PHE	CA-CB-CG	5.83	119.63	113.80
1	A	103	ASP	CA-CB-CG	5.81	118.41	112.60
1	A	172	GLY	N-CA-C	-5.76	99.52	113.18
1	A	63	LEU	CA-C-N	5.67	132.38	121.54
1	A	63	LEU	C-N-CA	5.67	132.38	121.54
1	A	173	TYR	N-CA-CB	-5.66	100.92	110.49
1	A	147	GLU	CG-CD-OE2	5.55	131.16	118.40
1	A	65	ASN	O-C-N	5.54	129.08	122.26
1	A	102	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	111	TYR	N-CA-CB	-5.37	101.41	110.49
1	A	48	ASP	CA-CB-CG	5.23	117.83	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	172	GLY	Peptide
1	A	174	CYS	Peptide

5.2 Too-close contacts

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	1418	0	1335	15	0
2	A	21	0	30	0	0
3	A	172	0	0	8	0
All	All	1611	0	1365	15	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 5.

All (15) 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:147:GLU:HB3	3:A:221:HOH:O	1.94	0.67
1:A:158:GLU:HG3	3:A:247:HOH:O	1.93	0.66
1:A:2:ARG:NH1	1:A:158:GLU:HA	2.17	0.60
1:A:122:LEU:N	3:A:209:HOH:O	2.38	0.54
1:A:174:CYS:O	1:A:175:ASP:HB2	2.08	0.52
1:A:39:ASP:HA	1:A:168:ILE:CG2	2.40	0.52
1:A:147:GLU:HG3	3:A:303:HOH:O	2.10	0.51
1:A:141:PRO:O	1:A:142:SER:OG	2.26	0.51
1:A:131:ASP:N	3:A:209:HOH:O	2.39	0.51
1:A:40:ASN:OD1	1:A:172:GLY:O	2.29	0.49
1:A:166:ARG:NH1	3:A:339:HOH:O	2.49	0.46
1:A:150:LYS:NZ	3:A:268:HOH:O	2.49	0.45
1:A:62:ARG:HD3	1:A:66:ASN:OD1	2.20	0.42
1:A:117:GLN:N	3:A:240:HOH:O	2.54	0.41
1:A:141:PRO:C	1:A:142:SER:HG	2.27	0.41

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	173/183 (94%)	164 (95%)	5 (3%)	4 (2%)	5 0

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	LEU
1	A	142	SER
1	A	173	TYR
1	A	111	TYR

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	151/158 (96%)	140 (93%)	11 (7%)	13 1

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	49	GLU
1	A	64	LEU
1	A	67	TRP
1	A	87	LYS
1	A	147	GLU
1	A	150	LYS
1	A	153	ARG
1	A	158	GLU
1	A	167	LEU
1	A	175	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	149	GLN
1	A	156	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 oligosaccharides 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	RTL	A	184	-	21,21,21	1.83	5 (23%)	28,28,28	1.94	7 (25%)

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	RTL	A	184	-	-	1/14/31/31	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	184	RTL	C5-C6	5.39	1.43	1.34
2	A	184	RTL	C20-C13	3.29	1.57	1.50
2	A	184	RTL	C19-C9	3.05	1.57	1.50
2	A	184	RTL	C8-C9	-2.37	1.40	1.46
2	A	184	RTL	C14-C13	2.11	1.38	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	184	RTL	C15-C14-C13	4.86	134.50	126.61
2	A	184	RTL	C11-C10-C9	3.57	132.28	127.28
2	A	184	RTL	C18-C5-C6	3.09	127.86	124.48
2	A	184	RTL	C17-C1-C6	-3.09	105.40	110.24
2	A	184	RTL	C7-C8-C9	2.83	130.42	126.23
2	A	184	RTL	C4-C5-C6	-2.63	119.15	122.70
2	A	184	RTL	C3-C4-C5	2.63	118.74	114.06

There are no chirality outliers.

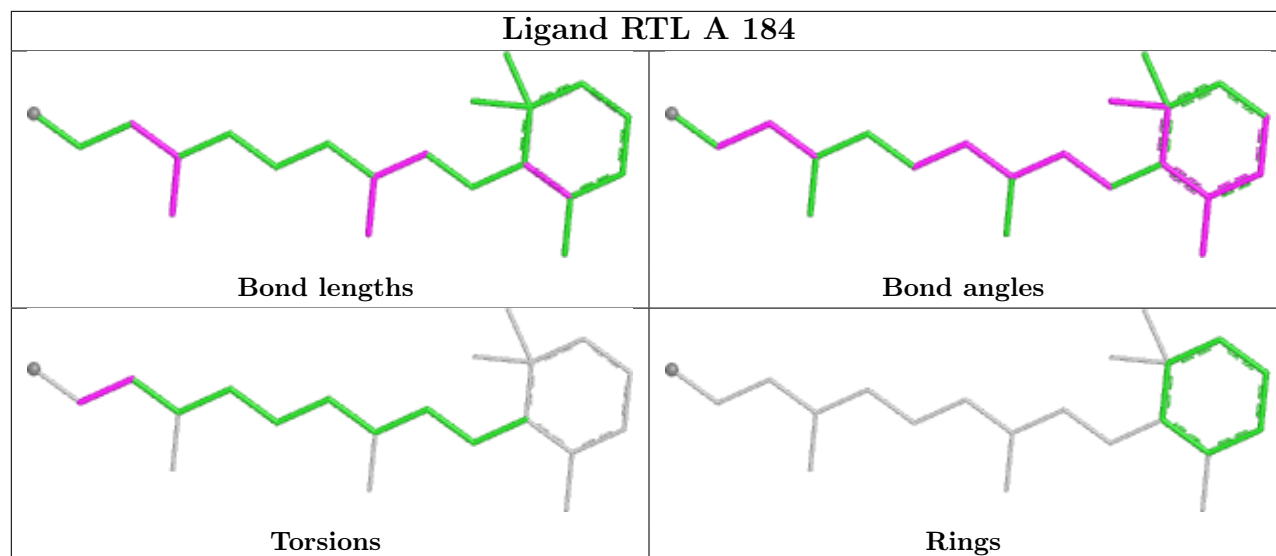
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	184	RTL	C13-C14-C15-O1

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	172:GLY	C	173:TYR	N	1.17

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.