



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 08:24 pm GMT

PDB ID : 1AD6
Title : DOMAIN A OF HUMAN RETINOBLASTOMA TUMOR SUPPRESSOR
Authors : Kim, H.Y.; Cho, Y.
Deposited on : 1997-02-21
Resolution : 2.30 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
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

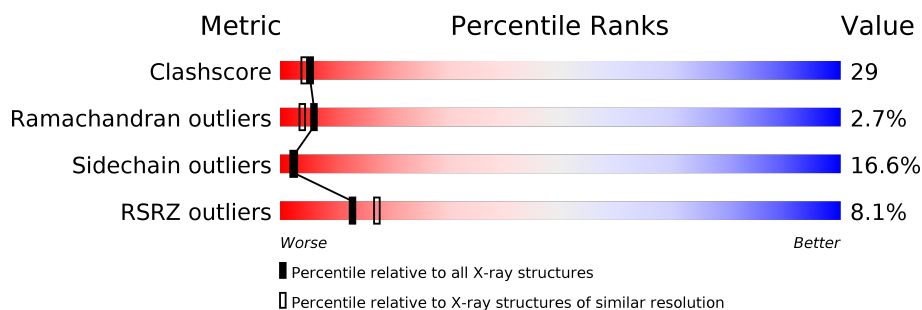
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	185	<div> <div>8%</div> <div>50%</div> <div>37%</div> <div>12%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1567 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 RETINOBLASTOMA TUMOR SUPPRESSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1491	947	252	279	13			

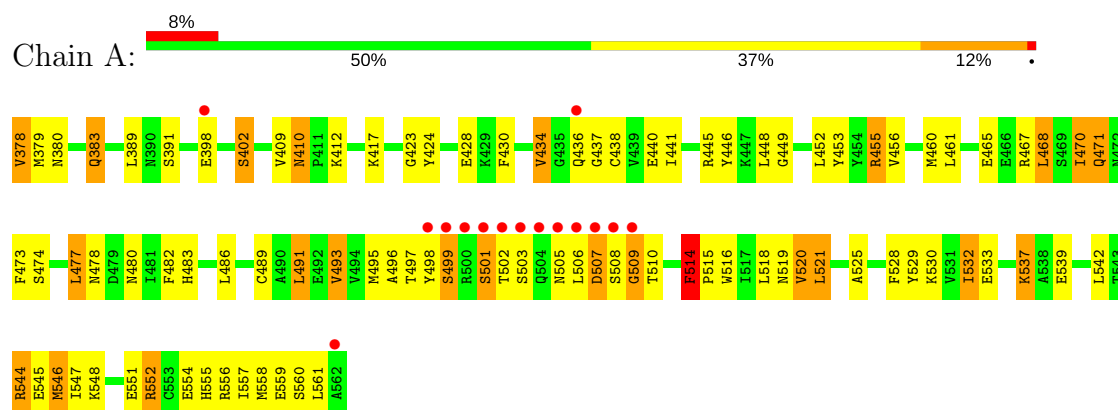
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	76	Total	O	0	0
			76	76		

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 the various outlier classes 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 ($\text{RSRZ} > 2$). 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: RETINOBLASTOMA TUMOR SUPPRESSOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	38.60Å 68.04Å 80.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.30 19.37 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.0 (6.00-2.30) 93.9 (19.37-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.38 (at 2.30Å)	Xtriage
Refinement program	TNT V. 5-D	Depositor
R, R_{free}	0.201 , 0.240 0.207 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 99.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	1567	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% 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](#)

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	0/1514	0.86	4/2032 (0.2%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	455	ARG	NE-CZ-NH2	-7.00	116.80	120.30
1	A	514	PHE	C-N-CD	-5.51	108.48	120.60
1	A	509	GLY	N-CA-C	-5.46	99.46	113.10
1	A	521	LEU	CA-CB-CG	5.41	127.75	115.30

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	1491	0	1512	87	0
2	A	76	0	0	6	0
All	All	1567	0	1512	87	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 29.

All (87) 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:468:LEU:HB3	1:A:470:ILE:HG23	1.51	0.93
1:A:503:SER:OG	1:A:506:LEU:HB2	1.80	0.81
1:A:441:ILE:HD11	1:A:507:ASP:HB2	1.65	0.78
1:A:449:GLY:HA3	1:A:491:LEU:HD13	1.67	0.76
1:A:441:ILE:CD1	1:A:507:ASP:HB2	2.18	0.73
1:A:465:GLU:HG3	1:A:470:ILE:O	1.88	0.73
1:A:529:TYR:CE2	1:A:554:GLU:HG2	2.24	0.73
1:A:502:THR:HG23	1:A:509:GLY:O	1.90	0.71
1:A:498:TYR:O	1:A:501:SER:N	2.24	0.71
1:A:514:PHE:HZ	1:A:557:ILE:HD11	1.56	0.69
1:A:514:PHE:CZ	1:A:557:ILE:HD11	2.28	0.69
1:A:498:TYR:HB3	1:A:502:THR:H	1.57	0.68
1:A:471:GLN:NE2	2:A:64:HOH:O	2.31	0.64
1:A:445:ARG:HH11	1:A:445:ARG:HG3	1.61	0.63
1:A:467:ARG:HD2	2:A:16:HOH:O	2.00	0.61
1:A:533:GLU:O	1:A:537:LYS:HD2	2.00	0.60
1:A:491:LEU:O	1:A:495:MET:HG2	2.02	0.60
1:A:449:GLY:HA3	1:A:491:LEU:CD1	2.33	0.58
1:A:498:TYR:HB2	1:A:501:SER:HA	1.87	0.57
1:A:518:LEU:HD11	1:A:528:PHE:CD2	2.41	0.56
1:A:519:ASN:ND2	2:A:54:HOH:O	2.36	0.56
1:A:436:GLN:NE2	1:A:436:GLN:HA	2.21	0.55
1:A:438:CYS:SG	1:A:507:ASP:HB3	2.46	0.55
1:A:497:THR:HG22	1:A:499:SER:HA	1.90	0.54
1:A:477:LEU:O	1:A:483:HIS:HE1	1.92	0.53
1:A:423:GLY:HA2	1:A:446:TYR:CE2	2.44	0.52
1:A:498:TYR:CB	1:A:502:THR:H	2.23	0.52
1:A:452:LEU:O	1:A:456:VAL:HG23	2.10	0.52
1:A:440:GLU:CD	1:A:440:GLU:H	2.14	0.51
1:A:378:VAL:HG23	1:A:379:MET:N	2.25	0.50
1:A:445:ARG:NH1	1:A:445:ARG:HG3	2.27	0.50
1:A:438:CYS:O	1:A:441:ILE:HB	2.11	0.50
1:A:460:MET:HE1	1:A:482:PHE:CE1	2.46	0.50
1:A:489:CYS:O	1:A:493:VAL:HG13	2.11	0.50
1:A:539:GLU:HG3	1:A:542:LEU:HD13	1.94	0.49
1:A:558:MET:CE	1:A:559:GLU:HG2	2.43	0.49
1:A:520:VAL:CG2	1:A:521:LEU:HD23	2.43	0.48
1:A:417:LYS:HD2	2:A:23:HOH:O	2.14	0.48
1:A:496:ALA:HB3	1:A:546:MET:CE	2.44	0.48
1:A:424:TYR:O	1:A:428:GLU:HG2	2.14	0.47
1:A:498:TYR:CB	1:A:501:SER:HA	2.43	0.47
1:A:460:MET:HE1	1:A:486:LEU:HD13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:ASN:HD22	1:A:410:ASN:C	2.17	0.47
1:A:455:ARG:HD3	1:A:539:GLU:OE1	2.14	0.47
1:A:410:ASN:ND2	1:A:412:LYS:H	2.12	0.47
1:A:460:MET:HE3	1:A:482:PHE:HZ	1.80	0.47
1:A:545:GLU:HG3	2:A:19:HOH:O	2.14	0.47
1:A:430:PHE:O	1:A:434:VAL:HB	2.15	0.47
1:A:498:TYR:O	1:A:501:SER:HA	2.14	0.47
1:A:515:PRO:HB2	1:A:519:ASN:HD21	1.79	0.46
1:A:468:LEU:HB3	1:A:470:ILE:CG2	2.35	0.46
1:A:468:LEU:CB	1:A:470:ILE:HG23	2.33	0.46
1:A:539:GLU:O	1:A:542:LEU:HB2	2.16	0.46
1:A:530:LYS:HB2	1:A:530:LYS:HE2	1.71	0.45
1:A:474:SER:O	1:A:478:ASN:OD1	2.34	0.45
1:A:430:PHE:HZ	1:A:508:SER:HB2	1.82	0.45
1:A:423:GLY:CA	1:A:446:TYR:CE2	3.00	0.45
1:A:525:ALA:HB1	1:A:557:ILE:HG12	1.99	0.45
1:A:453:TYR:OH	1:A:483:HIS:CD2	2.69	0.45
1:A:552:ARG:O	1:A:556:ARG:HG3	2.17	0.45
1:A:383:GLN:HB3	2:A:9:HOH:O	2.16	0.44
1:A:515:PRO:HB2	1:A:519:ASN:ND2	2.32	0.44
1:A:436:GLN:HE21	1:A:436:GLN:HA	1.81	0.44
1:A:460:MET:CE	1:A:482:PHE:CZ	3.00	0.44
1:A:470:ILE:HG12	1:A:473:PHE:CZ	2.53	0.44
1:A:502:THR:HG23	1:A:509:GLY:C	2.38	0.44
1:A:544:ARG:HG2	1:A:544:ARG:HH11	1.83	0.43
1:A:544:ARG:CG	1:A:544:ARG:HH11	2.31	0.43
1:A:552:ARG:NH1	1:A:555:HIS:ND1	2.66	0.43
1:A:430:PHE:CZ	1:A:441:ILE:HG22	2.53	0.43
1:A:423:GLY:HA2	1:A:446:TYR:CZ	2.54	0.43
1:A:410:ASN:HD22	1:A:412:LYS:H	1.67	0.42
1:A:507:ASP:N	1:A:507:ASP:OD1	2.52	0.42
1:A:516:TRP:O	1:A:520:VAL:HG13	2.19	0.42
1:A:547:ILE:O	1:A:551:GLU:HG3	2.19	0.42
1:A:496:ALA:HB3	1:A:546:MET:HE3	2.01	0.42
1:A:398:GLU:O	1:A:402:SER:OG	2.36	0.42
1:A:498:TYR:HB3	1:A:502:THR:N	2.28	0.42
1:A:441:ILE:HD12	1:A:507:ASP:HB2	2.01	0.42
1:A:520:VAL:HG22	1:A:521:LEU:HD23	2.03	0.41
1:A:507:ASP:O	1:A:508:SER:HB3	2.19	0.41
1:A:558:MET:O	1:A:561:LEU:HB2	2.21	0.41
1:A:456:VAL:O	1:A:460:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:SER:HG	1:A:506:LEU:HB2	1.82	0.41
1:A:532:ILE:HD12	1:A:554:GLU:HG3	2.03	0.41
1:A:503:SER:H	1:A:509:GLY:HA2	1.85	0.41
1:A:544:ARG:HD2	1:A:544:ARG:HA	1.85	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	183/185 (99%)	167 (91%)	11 (6%)	5 (3%)	6 4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	437	GLY
1	A	499	SER
1	A	501	SER
1	A	505	ASN
1	A	532	ILE

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	169/169 (100%)	141 (83%)	28 (17%)	2 2

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	VAL
1	A	380	ASN
1	A	383	GLN
1	A	389	LEU
1	A	391	SER
1	A	402	SER
1	A	409	VAL
1	A	410	ASN
1	A	434	VAL
1	A	448	LEU
1	A	461	LEU
1	A	468	LEU
1	A	470	ILE
1	A	471	GLN
1	A	477	LEU
1	A	480	ASN
1	A	491	LEU
1	A	493	VAL
1	A	507	ASP
1	A	510	THR
1	A	514	PHE
1	A	520	VAL
1	A	537	LYS
1	A	544	ARG
1	A	546	MET
1	A	548	LYS
1	A	552	ARG
1	A	560	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	384	GLN
1	A	390	ASN
1	A	399	ASN
1	A	405	ASN
1	A	410	ASN

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Mol	Chain	Res	Type
1	A	436	GLN
1	A	478	ASN
1	A	483	HIS
1	A	504	GLN
1	A	505	ASN
1	A	519	ASN
1	A	541	ASN

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](#)

There are no ligands in this entry.

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 [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, 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	185/185 (100%)	0.40	15 (8%) 13 17	11, 27, 67, 87	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	505	ASN	13.4
1	A	502	THR	11.1
1	A	500	ARG	10.1
1	A	501	SER	9.5
1	A	506	LEU	9.0
1	A	507	ASP	7.4
1	A	508	SER	7.2
1	A	503	SER	7.1
1	A	504	GLN	5.7
1	A	499	SER	5.7
1	A	562	ALA	5.6
1	A	509	GLY	3.4
1	A	498	TYR	2.7
1	A	436	GLN	2.3
1	A	398	GLU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.