



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2026 – 02:29 PM EDT

PDB ID : 37ZA / pdb_000037za
Title : Crystal structure of TP53 (89-293, Y220C)
Authors : Setser, J.W.
Deposited on : 2026-08-14
Resolution : 2.11 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
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

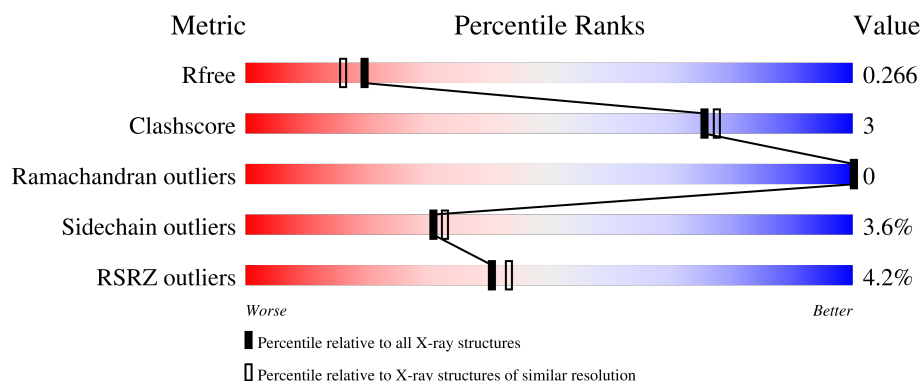
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

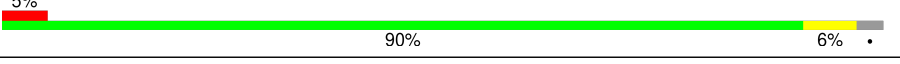
The reported resolution of this entry is 2.11 Å.

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}	180053	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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 $\leq 5\%$. 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	AAA	206	 3% 83% 14% ..
1	BBB	206	 5% 90% 6% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3297 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 Cellular tumor antigen p53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	201	Total	C	N	O	S	0	3	0
			1592	982	294	296	20			
1	BBB	200	Total	C	N	O	S	0	1	0
			1575	971	293	293	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	88	GLY	-	expression tag	UNP P04637
AAA	220	CYS	TYR	conflict	UNP P04637
BBB	88	GLY	-	expression tag	UNP P04637
BBB	220	CYS	TYR	conflict	UNP P04637

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	1	Total	Zn	0	0
			1	1		
2	BBB	1	Total	Zn	0	0
			1	1		

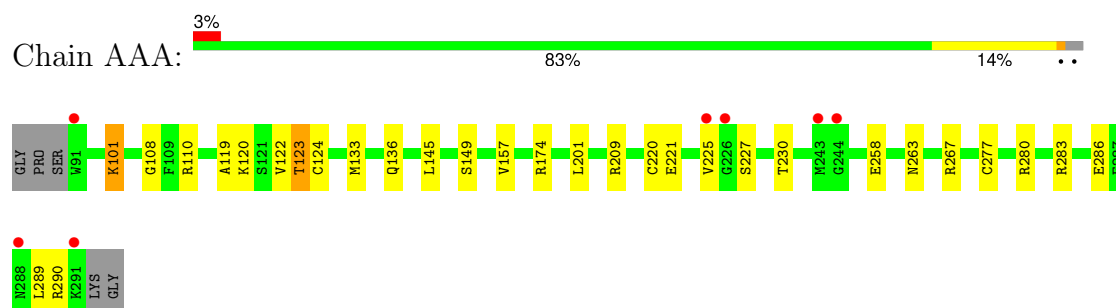
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	83	Total	O	0	0
			83	83		
3	BBB	45	Total	O	0	0
			45	45		

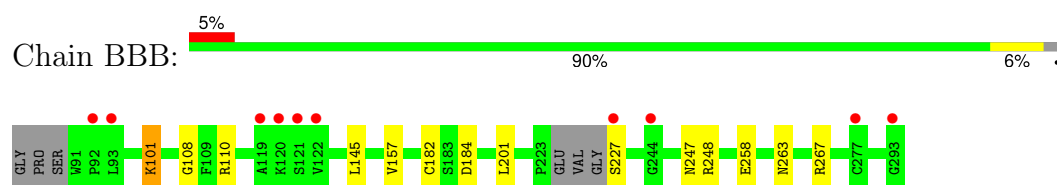
3 Residue-property plots [i](#)

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 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 ($RSRZ > 2$). 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: Cellular tumor antigen p53



- Molecule 1: Cellular tumor antigen p53



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.31Å 68.44Å 60.12Å 90.00° 98.81° 90.00°	Depositor
Resolution (Å)	44.87 – 2.11 44.87 – 2.11	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.87-2.11) 99.9 (44.87-2.11)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.202 , 0.263 0.208 , 0.266	Depositor DCC
R_{free} test set	1173 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.6	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	0.95	EDS
Total number of atoms	3297	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.78% 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: ZN

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	AAA	0.99	0/1635	1.24	0/2218
1	BBB	1.01	0/1614	1.23	0/2184
All	All	1.00	0/3249	1.23	0/4402

There are no bond length outliers.

There are no bond angle outliers.

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	AAA	1592	0	1559	12	0
1	BBB	1575	0	1542	7	0
2	AAA	1	0	0	0	0
2	BBB	1	0	0	0	0
3	AAA	83	0	0	1	0
3	BBB	45	0	0	0	0
All	All	3297	0	3101	19	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 3.

All (19) 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:AAA:124:CYS:SG	1:AAA:133[B]:MET:SD	2.68	0.91
1:AAA:277[B]:CYS:SG	1:AAA:280:ARG:NH2	2.67	0.67
1:AAA:123:THR:HA	1:AAA:136:GLN:NE2	2.13	0.64
1:AAA:221:GLU:O	1:AAA:230:THR:HG21	2.06	0.56
1:AAA:145:LEU:HD13	1:AAA:157:VAL:HG21	1.87	0.55
1:BBB:145:LEU:HD13	1:BBB:157:VAL:HG21	1.88	0.55
1:AAA:101:LYS:O	1:AAA:267:ARG:HD2	2.10	0.51
1:BBB:108:GLY:O	1:BBB:110:ARG:NH1	2.46	0.49
1:BBB:247:ASN:O	1:BBB:248:ARG:HB2	2.12	0.49
1:AAA:286:GLU:O	1:AAA:289:LEU:HB2	2.14	0.48
1:BBB:101:LYS:O	1:BBB:267:ARG:HD2	2.14	0.48
1:AAA:108:GLY:O	1:AAA:110:ARG:NH1	2.48	0.47
1:AAA:258:GLU:HA	1:AAA:263:ASN:O	2.17	0.43
1:AAA:220:CYS:O	3:AAA:401:HOH:O	2.21	0.43
1:BBB:258:GLU:HA	1:BBB:263:ASN:O	2.19	0.42
1:AAA:119:ALA:O	1:AAA:122:VAL:HG22	2.19	0.41
1:BBB:182[A]:CYS:SG	1:BBB:184:ASP:OD2	2.79	0.41
1:BBB:182[A]:CYS:SG	1:BBB:184:ASP:CG	3.04	0.41
1:AAA:201:LEU:HD23	1:AAA:201:LEU:HA	1.94	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	AAA	202/206 (98%)	200 (99%)	2 (1%)	0	100	100
1	BBB	197/206 (96%)	195 (99%)	2 (1%)	0	100	100
All	All	399/412 (97%)	395 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	AAA	184/184 (100%)	174 (95%)	10 (5%)	20	18
1	BBB	181/184 (98%)	178 (98%)	3 (2%)	53	61
All	All	365/368 (99%)	352 (96%)	13 (4%)	31	33

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	101	LYS
1	AAA	120	LYS
1	AAA	123	THR
1	AAA	149	SER
1	AAA	174	ARG
1	AAA	209	ARG
1	AAA	225	VAL
1	AAA	227	SER
1	AAA	283	ARG
1	AAA	290	ARG
1	BBB	101	LYS
1	BBB	201	LEU
1	BBB	227	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 [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	AAA	201/206 (97%)	0.33	7 (3%) 47 50	20, 38, 59, 84	3 (1%)
1	BBB	200/206 (97%)	0.55	10 (5%) 34 37	31, 45, 66, 94	1 (0%)
All	All	401/412 (97%)	0.44	17 (4%) 40 43	20, 41, 64, 94	4 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	244	GLY	4.4
1	AAA	226	GLY	4.1
1	BBB	92	PRO	3.4
1	AAA	225	VAL	3.2
1	BBB	119	ALA	2.5
1	AAA	291	LYS	2.5
1	BBB	122	VAL	2.4
1	BBB	244	GLY	2.3
1	AAA	243	MET	2.3
1	AAA	91	TRP	2.3
1	BBB	93	LEU	2.3
1	BBB	293	GLY	2.3
1	BBB	227	SER	2.2
1	AAA	288	ASN	2.2
1	BBB	121	SER	2.2
1	BBB	277	CYS	2.0
1	BBB	120	LYS	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	AAA	301	1/1	0.99	0.03	42,42,42,42	0
2	ZN	BBB	301	1/1	1.00	0.01	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.