



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:53 PM UTC

PDB ID : 6FF9 / pdb_00006ff9
Title : Mutant R280K of human P53
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Deposited on : 2018-01-04
Resolution : 2.00 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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

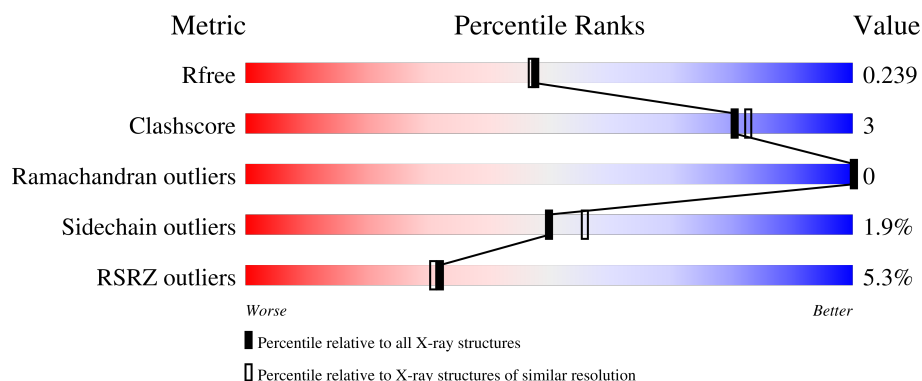
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	A	193	5% 94% 6%
1	B	193	7% 91% 8%
1	C	193	5% 92% 8%
1	D	193	4% 90% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6413 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	A	193	Total	C	N	O	S	0	1	0
			1517	938	278	284	17			
1	B	193	Total	C	N	O	S	0	0	0
			1515	936	278	285	16			
1	C	193	Total	C	N	O	S	0	2	0
			1526	943	279	286	18			
1	D	193	Total	C	N	O	S	0	0	0
			1515	936	278	285	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	280	LYS	ARG	engineered mutation	UNP P04637
B	280	LYS	ARG	engineered mutation	UNP P04637
C	280	LYS	ARG	engineered mutation	UNP P04637
D	280	LYS	ARG	engineered mutation	UNP P04637

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

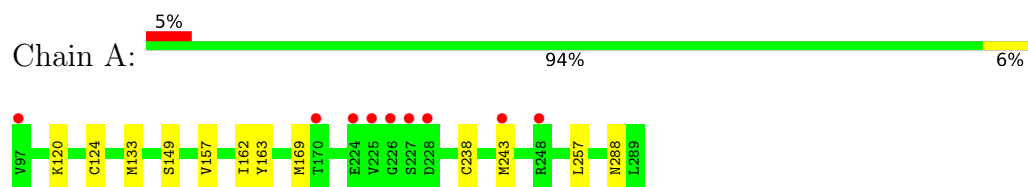
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	84	Total 84	O 84	0	0
3	B	85	Total 85	O 85	0	0
3	C	83	Total 83	O 83	0	0
3	D	84	Total 84	O 84	0	0

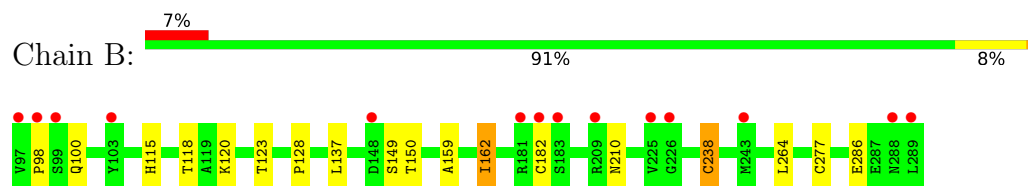
3 Residue-property plots [i](#)

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.

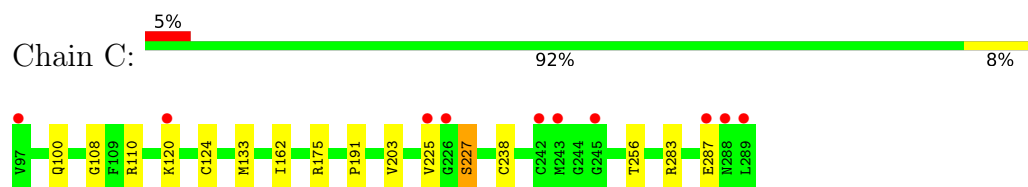
- Molecule 1: Cellular tumor antigen p53



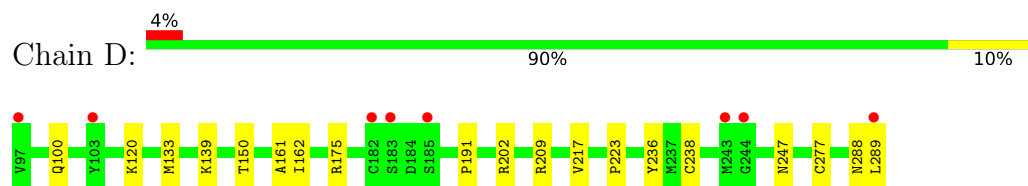
- Molecule 1: Cellular tumor antigen p53



- 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	68.57Å 69.44Å 83.33Å 90.00° 90.04° 90.00°	Depositor
Resolution (Å)	41.67 – 2.00 41.67 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.8 (41.67-2.00) 97.3 (41.67-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.00Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.182 , 0.226 0.196 , 0.239	Depositor DCC
R_{free} test set	2516 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -k,-h,-l 0.001 for k,h,-l 0.099 for h,-k,-l	Xtriage
Reported twinning fraction	0.751 for H, K, L 0.249 for -h,-k,l	Depositor
Outliers	0 of 51689 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6413	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 14.41% 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

5.1 Standard geometry

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	A	1.27	0/1555	1.11	0/2109
1	B	1.21	2/1550 (0.1%)	1.12	3/2102 (0.1%)
1	C	1.26	2/1564 (0.1%)	1.13	1/2121 (0.0%)
1	D	1.28	4/1550 (0.3%)	1.15	4/2102 (0.2%)
All	All	1.26	8/6219 (0.1%)	1.13	8/8434 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	118	THR	C-O	-5.71	1.15	1.24
1	C	203	VAL	CA-CB	5.71	1.60	1.54
1	C	162	ILE	N-CA	5.55	1.52	1.46
1	B	159	ALA	N-CA	-5.50	1.39	1.46
1	D	161	ALA	C-O	-5.38	1.17	1.23
1	D	139	LYS	C-O	5.25	1.30	1.23
1	D	162	ILE	N-CA	5.23	1.52	1.46
1	D	217	VAL	CA-C	5.04	1.58	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	223	PRO	N-CA-C	7.51	123.02	111.15
1	D	150	THR	CA-C-N	-6.51	114.91	119.66
1	D	150	THR	C-N-CA	-6.51	114.91	119.66
1	B	162	ILE	CB-CA-C	-5.86	100.22	110.71
1	C	227	SER	N-CA-CB	5.56	119.41	110.85
1	D	288	ASN	N-CA-C	5.24	118.77	112.38
1	B	150	THR	CA-C-N	-5.01	115.22	120.38
1	B	150	THR	C-N-CA	-5.01	115.22	120.38

There are no chirality outliers.

There are no planarity outliers.

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	1517	0	1480	7	0
1	B	1515	0	1475	12	0
1	C	1526	0	1488	8	0
1	D	1515	0	1475	10	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	84	0	0	0	0
3	B	85	0	0	4	0
3	C	83	0	0	3	0
3	D	84	0	0	2	0
All	All	6413	0	5918	31	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 (31) 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:D:133:MET:HE1	1:D:236:TYR:CE1	2.05	0.91
1:C:124:CYS:SG	1:C:133[B]:MET:SD	2.86	0.73
1:B:120:LYS:NZ	1:B:277:CYS:SG	2.65	0.70
1:C:227:SER:OG	3:C:401:HOH:O	2.17	0.58
1:B:115:HIS:HD2	3:B:481:HOH:O	1.86	0.58
1:D:202:ARG:HD3	3:D:401:HOH:O	2.03	0.58
1:A:243:MET:CE	1:B:264:LEU:HD23	2.35	0.57
1:A:124:CYS:SG	1:A:133:MET:SD	3.03	0.57
1:C:256:THR:HG21	3:C:461:HOH:O	2.03	0.56
1:D:133:MET:CE	1:D:236:TYR:CE1	2.84	0.55
1:D:133:MET:HE1	1:D:236:TYR:HE1	1.65	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:THR:OG1	3:B:401:HOH:O	2.19	0.53
1:A:163:TYR:O	1:A:169:MET:HG2	2.09	0.53
1:D:202:ARG:NH1	3:D:401:HOH:O	2.31	0.53
1:B:98:PRO:HG2	1:B:162:ILE:HD11	1.90	0.52
1:B:210:ASN:HA	1:D:247:ASN:HD22	1.74	0.52
1:C:283:ARG:O	1:C:287:GLU:HG2	2.13	0.48
1:C:225:VAL:HG11	1:D:100:GLN:NE2	2.30	0.46
1:D:120:LYS:NZ	1:D:277:CYS:SG	2.74	0.45
1:A:243:MET:HE1	1:B:264:LEU:HD23	2.00	0.44
1:A:162:ILE:HD12	1:A:169:MET:HE3	2.00	0.44
1:A:243:MET:HE1	1:B:264:LEU:CD2	2.48	0.43
1:B:137:LEU:HD12	1:B:238:CYS:C	2.45	0.42
1:C:175:ARG:HD3	1:C:191:PRO:O	2.20	0.42
1:D:175:ARG:HD3	1:D:191:PRO:O	2.19	0.41
1:A:157:VAL:HG22	1:A:257:LEU:HD23	2.02	0.41
1:C:108:GLY:O	1:C:110:ARG:NH1	2.54	0.41
1:B:128:PRO:HD2	3:B:467:HOH:O	2.20	0.41
1:C:100:GLN:HG2	3:C:409:HOH:O	2.21	0.41
1:B:210:ASN:HA	1:D:247:ASN:ND2	2.36	0.40
1:B:286:GLU:OE1	3:B:402:HOH:O	2.22	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	A	192/193 (100%)	192 (100%)	0	0	100	100
1	B	191/193 (99%)	191 (100%)	0	0	100	100
1	C	193/193 (100%)	192 (100%)	1 (0%)	0	100	100
1	D	191/193 (99%)	190 (100%)	1 (0%)	0	100	100
All	All	767/772 (99%)	765 (100%)	2 (0%)	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	A	174/173 (101%)	170 (98%)	4 (2%)	44	49
1	B	173/173 (100%)	169 (98%)	4 (2%)	44	49
1	C	175/173 (101%)	173 (99%)	2 (1%)	65	73
1	D	173/173 (100%)	170 (98%)	3 (2%)	53	60
All	All	695/692 (100%)	682 (98%)	13 (2%)	50	56

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

Mol	Chain	Res	Type
1	A	120	LYS
1	A	149	SER
1	A	238	CYS
1	A	288	ASN
1	B	100	GLN
1	B	149	SER
1	B	182	CYS
1	B	238	CYS
1	C	120	LYS
1	C	238	CYS
1	D	209	ARG
1	D	238	CYS
1	D	289	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	HIS
1	A	179	HIS
1	A	288	ASN
1	B	179	HIS

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Mol	Chain	Res	Type
1	C	100	GLN
1	C	179	HIS
1	C	210	ASN
1	D	115	HIS
1	D	179	HIS
1	D	210	ASN
1	D	247	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 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 ⓘ

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	193/193 (100%)	0.51	9 (4%)	36	35	15, 29, 48, 68	1 (0%)
1	B	193/193 (100%)	0.59	14 (7%)	21	19	18, 30, 54, 71	0
1	C	193/193 (100%)	0.46	10 (5%)	33	31	12, 28, 47, 70	2 (1%)
1	D	193/193 (100%)	0.51	8 (4%)	41	40	18, 28, 49, 62	0
All	All	772/772 (100%)	0.52	41 (5%)	32	31	12, 29, 50, 71	3 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	289	LEU	4.4
1	A	225	VAL	4.3
1	B	289	LEU	4.2
1	D	289	LEU	3.9
1	B	243	MET	3.9
1	C	97	VAL	3.6
1	B	183	SER	3.6
1	B	98	PRO	3.4
1	B	103	TYR	3.4
1	D	243	MET	3.4
1	C	225	VAL	3.1
1	D	183	SER	3.0
1	A	226	GLY	3.0
1	C	226	GLY	2.9
1	C	288	ASN	2.6
1	A	228	ASP	2.6
1	D	97	VAL	2.5
1	A	170	THR	2.5
1	B	226	GLY	2.5
1	C	120	LYS	2.5
1	B	99	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	97	VAL	2.4
1	A	243	MET	2.4
1	C	287	GLU	2.3
1	B	288	ASN	2.3
1	D	185	SER	2.3
1	B	225	VAL	2.2
1	A	224	GLU	2.2
1	D	244	GLY	2.2
1	B	182	CYS	2.2
1	D	182	CYS	2.2
1	D	103	TYR	2.2
1	B	97	VAL	2.1
1	B	209	ARG	2.1
1	C	245	GLY	2.1
1	C	243	MET	2.1
1	A	227	SER	2.1
1	C	242	CYS	2.0
1	A	248	ARG	2.0
1	B	181	ARG	2.0
1	B	148	ASP	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(Å ²)	Q<0.9
2	ZN	D	301	1/1	0.98	0.03	30,30,30,30	0
2	ZN	C	301	1/1	0.99	0.04	24,24,24,24	0
2	ZN	B	301	1/1	0.99	0.02	32,32,32,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	301	1/1	1.00	0.05	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.