



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 08:24 pm BST

PDB ID : 4CRI  
Title : Crystal Structure of 53BP1 tandem tudor domains in complex with methylated K810 Rb peptide  
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Deposited on : 2014-02-26  
Resolution : 2.35 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

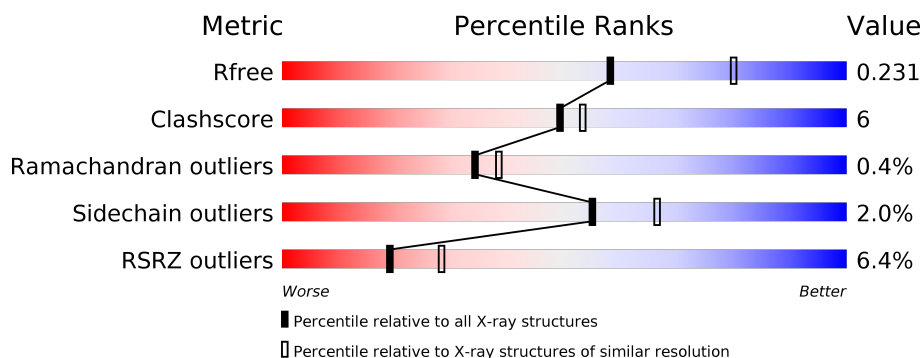
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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))
$R_{free}$	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 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	176	<div> <div>4%</div> <div> <div></div> <div>61%</div> <div>7%</div> <div>32%</div> </div> </div>
1	B	176	<div> <div>5%</div> <div> <div></div> <div>55%</div> <div>14%</div> <div>32%</div> </div> </div>
2	C	17	<div> <div>6%</div> <div> <div>18%</div> <div>24%</div> <div>59%</div> </div> </div>
2	D	17	<div> <div>24%</div> <div>12%</div> <div>65%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1991 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 TUMOR SUPPRESSOR P53-BINDING PROTEIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	S	0	0	1
			944	605	157	179	3			
1	B	120	Total	C	N	O	S	0	0	1
			922	589	152	178	3			

- Molecule 2 is a protein called RB1 PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	7	Total	C	N	O	0	0	0
			60	43	8	9			
2	D	6	Total	C	N	O	0	0	0
			42	28	7	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	818	CYS	-	expression tag	UNP P78495
D	818	CYS	-	expression tag	UNP P78495

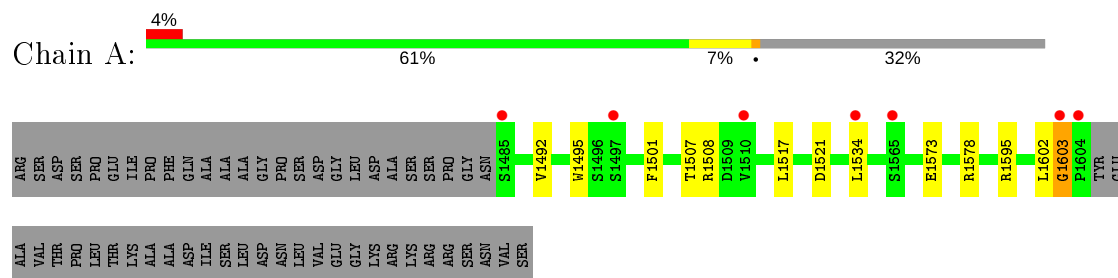
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		
3	B	6	Total	O	0	0
			6	6		

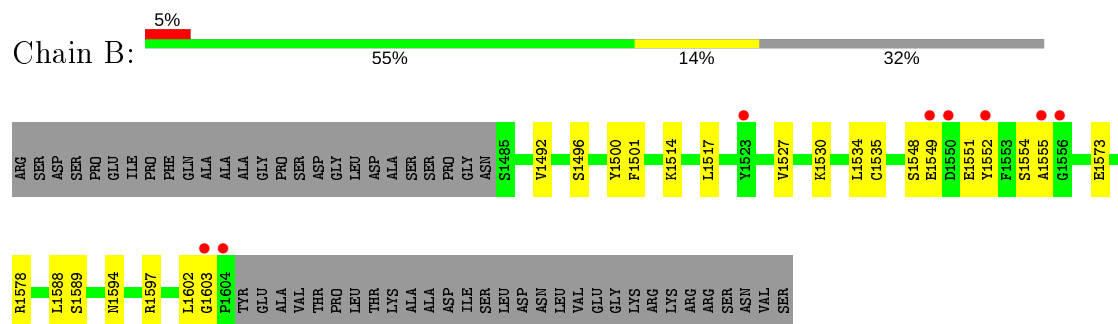
### 3 Residue-property plots [i](#)

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 ( $RSRZ > 2$ ). 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.

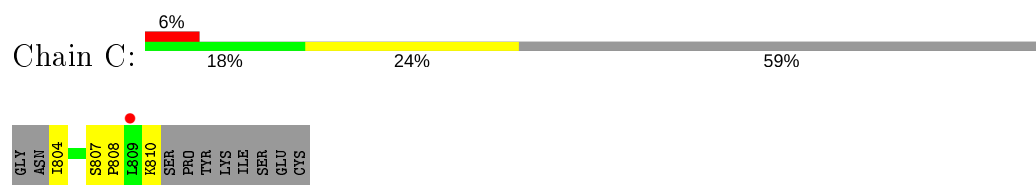
#### • Molecule 1: TUMOR SUPPRESSOR P53-BINDING PROTEIN 1



#### • Molecule 1: TUMOR SUPPRESSOR P53-BINDING PROTEIN 1



#### • Molecule 2: RB1 PROTEIN



#### • Molecule 2: RB1 PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	105.79Å 105.79Å 156.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.51 – 2.35 39.51 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.51-2.35) 99.9 (39.51-2.35)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 2.34Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.194 , 0.230 0.199 , 0.231	Depositor DCC
$R_{free}$ test set	1160 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	59.6	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 47.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	1991	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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: MLY

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	0.51	1/963 (0.1%)	0.65	1/1295 (0.1%)
1	B	0.47	1/941 (0.1%)	0.67	0/1272
2	C	0.57	0/50	0.73	0/68
2	D	0.35	0/31	0.64	0/42
All	All	0.50	2/1985 (0.1%)	0.66	1/2677 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1603	GLY	C-N	-5.55	1.23	1.34
1	A	1603	GLY	C-N	-5.16	1.24	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1517	LEU	CB-CG-CD2	-5.88	101.00	111.00

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	944	0	925	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	922	0	871	14	0
2	C	60	0	68	4	0
2	D	42	0	41	4	0
3	A	17	0	0	1	1
3	B	6	0	0	0	0
All	All	1991	0	1905	23	1

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 6.

All (23) 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:B:1594:ASN:OD1	1:B:1597:ARG:NH2	2.15	0.79
1:B:1573:GLU:HB2	1:B:1578:ARG:HG2	1.64	0.77
1:B:1549:GLU:HA	2:D:805:TYR:HA	1.65	0.76
1:A:1573:GLU:HB2	1:A:1578:ARG:HG2	1.78	0.65
1:B:1552:TYR:HA	2:D:806:ILE:HG12	1.86	0.58
1:B:1548:SER:H	2:D:806:ILE:HG22	1.70	0.57
1:A:1595:ARG:NH1	3:A:2017:HOH:O	2.33	0.56
1:B:1492:VAL:HG12	1:B:1501:PHE:HB3	1.88	0.54
1:A:1521:ASP:OD2	2:C:810:MLY:HH22	2.08	0.53
1:B:1534:LEU:HD22	1:B:1602:LEU:HD11	1.93	0.51
1:B:1514:LYS:NZ	2:C:810:MLY:O	2.44	0.50
1:A:1507:THR:O	1:A:1508:ARG:HG2	2.11	0.50
1:B:1554:SER:OG	1:B:1555:ALA:N	2.49	0.45
1:B:1548:SER:HB3	1:B:1551:GLU:O	2.19	0.43
2:C:807:SER:HB2	2:C:808:PRO:HD2	2.01	0.43
1:A:1495:TRP:CD1	2:C:810:MLY:HD2	2.54	0.43
1:B:1500:TYR:CD2	1:B:1589:SER:HA	2.55	0.42
1:A:1492:VAL:HG12	1:A:1501:PHE:HB3	2.02	0.42
1:B:1517:LEU:HD13	1:B:1527:VAL:HB	2.02	0.42
1:B:1551:GLU:O	2:D:806:ILE:HG21	2.20	0.41
1:A:1534:LEU:HD22	1:A:1602:LEU:HD11	2.02	0.41
1:A:1534:LEU:HB2	1:A:1602:LEU:CD1	2.50	0.41
1:B:1501:PHE:HB2	1:B:1588:LEU:HB2	2.01	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:2015:HOH:O	3:A:2015:HOH:O[12_565]	1.63	0.57

## 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	118/176 (67%)	115 (98%)	2 (2%)	1 (1%)	19	20
1	B	118/176 (67%)	114 (97%)	4 (3%)	0	100	100
2	C	5/17 (29%)	5 (100%)	0	0	100	100
2	D	4/17 (24%)	4 (100%)	0	0	100	100
All	All	245/386 (64%)	238 (97%)	6 (2%)	1 (0%)	34	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1603	GLY

### 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	98/146 (67%)	98 (100%)	0	100	100
1	B	93/146 (64%)	90 (97%)	3 (3%)	39	47
2	C	6/15 (40%)	5 (83%)	1 (17%)	2	1
2	D	3/15 (20%)	3 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	200/322 (62%)	196 (98%)	4 (2%)	55 66

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1496	SER
1	B	1530	LYS
1	B	1535	CYS
2	C	804	ILE

Some 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 ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
2	MLY	D	810	2	9,10,11	0.69	0	6,11,13	0.94	0
2	MLY	C	810	2	9,10,11	0.63	0	6,11,13	0.88	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
2	MLY	D	810	2	-	2/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLY	C	810	2	-	6/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	810	MLY	N-CA-CB-CG
2	C	810	MLY	C-CA-CB-CG
2	C	810	MLY	O-C-CA-CB
2	C	810	MLY	CG-CD-CE-NZ
2	D	810	MLY	CD-CE-NZ-CH1
2	C	810	MLY	CA-CB-CG-CD
2	D	810	MLY	CE-CD-CG-CB
2	C	810	MLY	CE-CD-CG-CB

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	810	MLY	3	0

## 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

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(Å²)	Q<0.9
1	A	120/176 (68%)	0.22	7 (5%)	23 33	42, 66, 98, 122	0
1	B	120/176 (68%)	0.36	8 (6%)	17 26	52, 77, 126, 145	0
2	C	6/17 (35%)	1.05	1 (16%)	1 2	47, 51, 63, 81	0
2	D	5/17 (29%)	0.63	0	100 100	92, 96, 115, 130	0
All	All	251/386 (65%)	0.31	16 (6%)	19 28	42, 71, 115, 145	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1604	PRO	13.3
1	B	1552	TYR	5.1
1	A	1604	PRO	4.7
1	B	1556	GLY	3.8
1	A	1510	VAL	3.6
1	B	1549	GLU	3.1
1	B	1555	ALA	3.0
1	A	1485	SER	2.7
2	C	809	LEU	2.4
1	B	1603	GLY	2.4
1	B	1550	ASP	2.3
1	B	1523	TYR	2.2
1	A	1565	SER	2.2
1	A	1603	GLY	2.1
1	A	1534	LEU	2.1
1	A	1497	SER	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	MLY	D	810	11/12	0.84	0.21	65,79,85,85	0
2	MLY	C	810	11/12	0.95	0.27	49,55,64,66	0

### 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.